

Versar^{inc.}
ENVIRONMENTAL RISK MANAGEMENT

Versar^{inc.}

ORIGINAL
(Red)

441596

**VOLATILE ANALYSIS
ANALYTICAL DATA PACKAGE**

**CLIENT : VERSAR DIV. 31
SITE : CDM
PROJECT: 420.1 B#2
CONTROL: 2536
DATE : 05/23/90**

DELIVERABLES INDEX -- VOA ANALYSES ONLY**I. CASE NARRATIVE**

- A. The narrative contains: case and contract numbers, summary of QC analyses, discussion of any analytical problems, and documentation of all corrective actions.
- B. Copy of Sample Traffic Reports.

II. QC SUMMARY

- A. Surrogate Percent Recovery Summary (Form II)
- B. Matrix Spike/Matrix Spike Duplicate Summary (Form III)
- C. Method Blank Summary (Form IV)
- D. GC/MS Tuning and Calibration Standard (Form V)

III. SAMPLE DATA

- A. Organic Analysis Data Sheet (Form I)
- B. Tentatively Identified Compounds (TIC)
- C. Raw Data for the Volatile Sample Fraction
 - 1. Reconstructed Ion Chromatogram (s)
 - 2. Quantitation Report
 - 3. Raw HSL mass spectra and background subtract HSL mass spectra with lab generated HSL standard spectra
 - 4. GC/MS library search spectra for each TIC

Note: Samples arranged in alpha-numeric sequence.

IV. STANDARDS DATA

- A. Initial Calibration Data (Form VI) in order if more than one instrument is used.
- B. VOA standard(s) reconstructed ion chromatograms and quantitation reports for the initial calibration(s).
- C. Continuing Calibration (Form VII) in order by instrument.
- D. VOA standard(s) reconstructed ion chromatograms and quantitation reports for the continuing calibration(s).
- E. Internal Standard Data (Form VIII).

V. RAW QC DATA

- A. BFB (For each 12-hour period, for each GC/MS system)
 - 1. Bar graph spectrum
 - 2. Mass listing

V. RAW QC DATA (CONT'D)**B. Reagent Blank Data**

1. Organic Analysis Data Sheet (Form I)
2. Tentatively Identified Compounds (TIC)
3. Raw Data
 - a. Reconstructed ion chromatogram(s) and quantitation report(s).
 - b. HSL spectra with lab generated standard
 - i. Raw HSL compound spectra
 - ii. Enhanced or background subtracted spectra
 - iii. Laboratory generated HSL standard spectra
 - iv. GC/MS library search spectra for each TIC
 - v. Quantitation/calculation of each TIC concentration

C. Matrix Spike/Matrix Spike Duplicate Data

1. Tabulated results (Form I) of non-spiked HSL compounds.
2. Reconstructed ion chromatograms and quantitation reports.

VI. SAMPLE PREPARATION

- A. Parameter Request Sheet
- B. Sample Receiving Log-In Information
- C. VOA Sample Comments
- D. Instrument Injection Logs
- E. Field Chain of Custody
- F. Copy(s) of Federal Express Airbills

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**ORIGINAL
(Red)**

I. NARRATIVE

II. QC SUMMARY

100003

May 22, 1990

Narrative

Versar Division 31 - CDM

Volatile Organic Analysis - EPA CLP Protocol

Versar Project 420.1 - Batch 2

Control 2536

This report contains the analytical data for the volatile organic analysis of six (6) sludge samples. The samples listed below arrived at Versar on April 19, 1990. Results from the Toxicity Characteristic Leaching Procedure (TCLP) will be reported separately. Results for additional analytical parameters, including semivolatile organic analysis and TCLP, are provided separately.

SAMPLE LIST

1	3	6
2	5	7

//////////

GC/MS instrument calibration using BFB met requirements for volatile organic analysis. SPCC and CCC criteria were met for the volatile initial calibration curve and the continuing calibration standards. All GC/MS analyses occurred during the twelve hour period that followed daily instrument calibration. All sample analyses for volatile organic compounds were completed within eight days of sample receipt.

Volatile surrogate standard recovery values and internal standard area abundances and associated relative retention times met specified QC limits for all samples except sample 1 and the QC analyses performed on this sample. Since the QC analyses and the sample analysis are replicates of one another, no reanalyses have been performed. One set of matrix spike and matrix spike duplicate (MS/MSD) QC analyses was analyzed for the soil matrix. All recovery and relative (RPD) values were within specified limits. The samples were analyzed without dilution. Nontarget compounds were tentatively identified using the EPA/NBS Mass Spectral Database Library.

Detection limits and concentrations of volatile organic compounds in the soil samples are reported on a "dry weight" basis, which accounts for the moisture content of the samples. The sludge samples were analyzed using protocol developed for the soil matrix and "SOIL" appears as the matrix on the reporting forms.

Definitions of data qualifier flags used on the individual data summary pages are provided in the listing which immediately follows this case narrative.

100004

Narrative - Page 2
Versar Div 31 - CDM
Control 2536

Please contact Janet Beckman, Laboratory Project Manager, should you have any questions or require additional information pertaining to the volatile organic analysis of these samples.

Data Release Approved By:



Linda E. Bock
GC/MS Data Quality Manager
Laboratory Operations

Narrative Reviewed By:



Debbie Maxwell
Lawrence P. Pollack
GC/MS Quality Assurance Manager

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Data Qualifier Flags

- J For Target Compounds: This flag is used when mass spectral data indicates the presence of a compound but the result is less than the specified detection limit but still greater than zero.
- For Non Target Compounds: This flag indicates that the concentration is an estimated value, assuming a 1 to 1 response with the internal standard.
- B This flag is used when the analyte is found in the blank as well as in the sample. It indicates possible/probable contamination and warns the data user to take appropriate action.
- u This flag states that the compound was analyzed for but was not detected. The number is the minimum attainable detection limit for the sample.
- X or T This flag states that the mass spectrum does not meet EPA CLP criteria for confirmation, but compound presence is strongly suspected.
- E This flag is used to indicate that the quantitation of the analyte is outside the linear calibration of the curve and that dilution was required in order to properly quantitate.
- D This flag is used to indicate the value for the target analyte was calculated from a dilution (see "E" flag above).
- Y This flag is used when a matrix spike compound is also confirmed present in the unspiked sample.

Flags excerpted from and established by the US EPA Contract Lab Program (CLP) protocol.

2B
SOIL VOLATILE SURROGATE RECOVERY

Lab Name: VERSAR INC.

Contract: _____

La. Code: VERSAR Case No.: R3-7

SAS No.: _____

SDG No.: 1 _____

Level: (low/med) LOW

EPA SAMPLE NO.	S1 (TOL) #	S2 (BFB) #	S3 (DCE) #	OTHER	TOT OUT
01 1	125 *	85	72	0	1
02 2	106	87	70	0	0
03 3	103	86	71	0	0
04 5	108	93	83	0	0
05 6	111	94	84	0	0
06 7	111	95	87	0	0
07 1MS	117	75	62 *	0	1
08 1MSD	117	77	62 *	0	1
09 VBLK87	97	93	87	0	0
10 VBLK12	102	100	84	0	0

QC LIMITS

S1 (TOL) = Toluene-d8 (81-117)

S2 (BFB) = Bromofluorobenzene (74-121)

S3 (DCE) = 1,2-Dichloroethane-d4 (70-121)

Column to be used to flag recovery values

* Values outside of contract required QC limits

D Surrogates diluted out

3B
SOIL VOLATILE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

ORIGINAL
(Rev)

Lab Name: VERSAR INC. Contract: _____

La Code: VERSAR Case No.: R3-7 SAS No.: _____ SDG No.: 1

Matrix Spike - EPA Sample No.: 1 Level: (low/med) LOW

COMPOUND	SPIKE ADDED (ug/Kg)	SAMPLE CONCENTRATION (ug/Kg)	MS CONCENTRATION (ug/Kg)	MS % REC #	QC LIMITS REC.
1,1-Dichloroethene	71.4	0	40.3	56 *	59-172
Trichloroethene	71.4	0	51.1	72	62-137
Benzene	71.4	0	78.6	110	66-142
Toluene	71.4	0	82.4	115	59-139
Chlorobenzene	71.4	0	67.0	94	60-133

COMPOUND	SPIKE ADDED (ug/Kg)	MSD CONCENTRATION (ug/Kg)	MSD % REC #	% RPD #	QC LIMITS RPD	REC.
1,1-Dichloroethene	71.4	41.1	58 *	-4	22	59-172
Trichloroethene	71.4	54.7	77	-7	24	62-137
Benzene	71.4	80.7	113	-3	21	66-142
Toluene	71.4	92.1	129	-11	21	59-139
Chlorobenzene	71.4	73.9	104	-10	21	60-133

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits

RPD: 0 out of 5 outside limits

Spike Recovery: 2 out of 10 outside limits

COMMENTS: CLP,,,VSTD20,L,W,17503,V,CC-020,25ML
INST R: SP1000 60/80 CARBOPACK 2M@45C > 225C@8C/M > 30M@225C

4A
VOLATILE METHOD BLANK SUMMARY

(PMT, 1981)
(Rev.)

Lab Name: VERSAR INC.

Contract: _____

Lab Code: VERSAR Case No.: R3-7 SAS No.: _____ SDG No.: 1 _____

Lab File ID: U3387 Lab Sample ID: VBLK87 _____

Date Analyzed: 04/26/90 Time Analyzed: 931 _____

Matrix: (soil/water) SOIL Level: (low/med) LOW _____

Instrument ID: U _____

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	TIME ANALYZED
01 2	16427	U3393	1431
02 3	16428	U3394	1524
03 1MS	16426MS	U3398	1811
04 1MSD	16426MSD	U3399	1853

COMMENTS: CLP,,,VBLK87,L,S,VBLK87,V,BLANK,,5 ML,,
INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

4A
VOLATILE METHOD BLANK SUMMARY

(M-A)

Lab Name: VERSAR INC.

Contract: _____

La. Code: VERSAR Case No.: R3-7 SAS No.: _____ SDG No.: 1 _____

Lab File ID: U3412 Lab Sample ID: VBLK12 _____

Date Analyzed: 04/27/90 Time Analyzed: 926 _____

Matrix: (soil/water) SOIL Level: (low/med) LOW _____

Instrument ID: U _____

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	TIME ANALYZED
01	1	16426	U3420	1614
02	5	16429	U3413	1022
03	6	16430	U3414	1105
04	7	16431	U3415	1147

COMMENTS: CLP,,,VBLK12,L,S,VBLK12,V,BLANK,,5 ML,,
INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

5A
 VOLATILE ORGANIC GC/MS TUNING AND MASS
 CALIBRATION - BROMOFLUOROBENZENE (BFB)

ON/ONLINE

(R)

Lab Name: VERSAR INC.

Contract: _____

Lab Code: VERSAR Case No.: R3-7

SAS No.: _____ SDG No.: 1 _____

Lab File ID: U3326

BFB Injection Date: 04/23/90

Instrument ID: U

BFB Injection Time: 1549

Matrix: (soil/water) SOIL Level: (low/med) LOW Column: (pack/cap) PACK

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	23.4
75	30.0 - 60.0% of mass 95	46.5
75	Base peak, 100% relative abundance	100.0
16	5.0 - 9.0% of mass 95	7.9
173	Less than 2.0% of mass 174	0.0 (0.0)1
174	Greater than 50.0% of mass 95	68.8
175	5.0 - 9.0% of mass 174	4.5 (6.6)1
176	Greater than 95.0%, but less than 101.0% of mass 174	66.9 (97.3)1
177	5.0 - 9.0% of mass 176	3.5 (5.3)2

1-Value is % mass 174

2-Value is % mass 176

THE TUNE APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01 VSTD050	22522	U3327	04/23/90	1623
02 VSTD200	22523	U3328	04/23/90	1721
03 VSTD150	22524	U3329	04/23/90	1809
04 VSTD100	22525	U3330	04/23/90	1852
05 VSTD020	22526	U3331	04/23/90	1935

5A
VOLATILE ORGANIC GC/MS TUNING AND MASS
CALIBRATION - BROMOFLUOROBENZENE (BFB)

Lab Name: VERSAR INC.

Contract: _____

Lab Code: VERSAR Case No.: R3-7

SAS No.: _____ SDG No.: 1 _____

Lab File ID: U3385

BFB Injection Date: 04/26/90

Instrument ID: U

BFB Injection Time: 0752

Matrix: (soil/water) SOIL Level: (low/med) LOW Column: (pack/cap) PACK

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	28.7
75	30.0 - 60.0% of mass 95	50.0
95	Base peak, 100% relative abundance	100.0
116	5.0 - 9.0% of mass 95	7.2
173	Less than 2.0% of mass 174	0.0 (0.0)1
174	Greater than 50.0% of mass 95	62.2
175	5.0 - 9.0% of mass 174	4.0 (6.5)1
176	Greater than 95.0%, but less than 101.0% of mass 174	61.6 (99.0)1
177	5.0 - 9.0% of mass 176	3.6 (5.8)2

1-Value is % mass 174

2-Value is % mass 176

THE TUNE APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01 VSTD050	22546	U3386	04/26/90	830
02 VBLK87	VBLK87	U3387	04/26/90	931
03 2	16427	U3393	04/26/90	1431
04 3	16428	U3394	04/26/90	1524
05 1MS	16426MS	U3398	04/26/90	1811
06 1MSD	16426MSD	U3399	04/26/90	1853

5A
VOLATILE ORGANIC GC/MS TUNING AND MASS
CALIBRATION - BROMOFLUOROBENZENE (BFB)

(R)
OVR

Lab Name: VERSAR INC.

Contract: _____

Lab Code: VERSAR Case No.: R3-7

SAS No.: _____ SDG No.: 1 _____

Lab File ID: U3410

BFB Injection Date: 04/27/90

Instrument ID: U

BFB Injection Time: 0745

Matrix: (soil/water) SOIL Level: (low/med) LOW Column: (pack/cap) PACK

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	28.0
75	30.0 - 60.0% of mass 95	52.4
95	Base peak, 100% relative abundance	100.0
16	5.0 - 9.0% of mass 95	7.2
173	Less than 2.0% of mass 174	0.0 (0.0)1
174	Greater than 50.0% of mass 95	73.1
175	5.0 - 9.0% of mass 174	5.5 (7.5)1
176	Greater than 95.0%, but less than 101.0% of mass 174	71.3 (97.6)1
177	5.0 - 9.0% of mass 176	4.4 (6.2)2

1-Value is % mass 174

2-Value is % mass 176

THE TUNE APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01 VSTD050	22554	U3411	04/27/90	821
02 VBLK12	VBLK12	U3412	04/27/90	926
03 5	16429	U3413	04/27/90	1022
04 6	16430	U3414	04/27/90	1105
05 7	16431	U3415	04/27/90	1147
06 1	16426	U3420	04/27/90	1614

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III. SAMPLE DATA PACKAGE

100014

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: VERSAR INC.

Contract: _____

ON/OWNL

1 (P.M.)

Lab Code: VERSAR Case No.: R3-7

SAS No.: _____

SDG No.: 1

Matrix: (soil/water) SOIL

Lab Sample ID: 16426

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: U3420

Level: (low/med) LOW

Date Received: 04/19/90

% Moisture: not dec. 30

Date Analyzed: 04/27/90

Column: (pack/cap) PACK

Dilution Factor: 1.0

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

74-87-3-----Chloromethane	14	U
74-83-9-----Bromomethane	14	U
75-01-4-----Vinyl chloride	14	U
75-00-3-----Chloroethane	14	U
75-09-2-----Methylene chloride	7	U
67-64-1-----Acetone	160	
75-15-0-----Carbon disulfide	7	U
75-35-4-----1,1-Dichloroethene	7	U
75-34-3-----1,1-Dichloroethane	7	U
540-59-0-----1,2-Dichloroethene (total)	7	U
67-66-3-----Chloroform	7	U
107-06-2-----1,2-Dichloroethane	7	U
78-93-3-----2-Butanone	14	U
71-55-6-----1,1,1-Trichloroethane	7	U
56-23-5-----Carbon tetrachloride	7	U
108-05-4-----Vinyl acetate	14	U
75-27-4-----Bromodichloromethane	7	U
78-87-5-----1,2-Dichloropropane	7	U
10061-01-5-----cis-1,3-Dichloropropene	7	U
79-01-6-----Trichloroethene	7	U
124-48-1-----Dibromochloromethane	7	U
79-00-5-----1,1,2-Trichloroethane	7	U
71-43-2-----Benzene	7	U
10061-02-6-----Trans-1,3-dichloropropene	7	U
75-25-2-----Bromoform	7	U
108-10-1-----4-Methyl-2-pentanone	14	U
591-78-6-----2-Hexanone	14	U
127-18-4-----Tetrachloroethene	7	U
79-34-5-----1,1,2,2-Tetrachloroethane	7	U
108-88-3-----Toluene	7	U
108-90-7-----Chlorobenzene	7	U
100-41-4-----Ethylbenzene	7	U
100-42-5-----Styrene	7	U
1330-20-7-----Total xylenes	7	U

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: VERSAR INC.

Contract: _____

1	ORIGINAL (REC)
---	-------------------

Lab Code: VERSAR Case No.: R3-7 SAS No.: _____ SDG No.: 1 _____

Matrix: (soil/water) SOIL

Lab Sample ID: 16426 _____

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: U3420 _____

Level: (low/med) LOW

Date Received: 04/19/90

* Moisture: not dec. 30

Date Analyzed: 04/27/90

Column (pack/cap) PACK

Dilution Factor: 1.0

No. 'ber TICs found: 4

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 106-97-8	BUTANE	8.45	9.7	J
2. 109-66-0	PENTANE (ACN) (DOT)	14.29	15	J
3. 110-54-3	HEXANE (DOT)	19.70	13	J
4.	UNKNOWN PAH	32.96	80	J

OPTIONAL
DATA

RIC
04/27/90 16:14:00

DATA: U3420 #1

SCANS 1 TO 840

CALI: U3420 #2

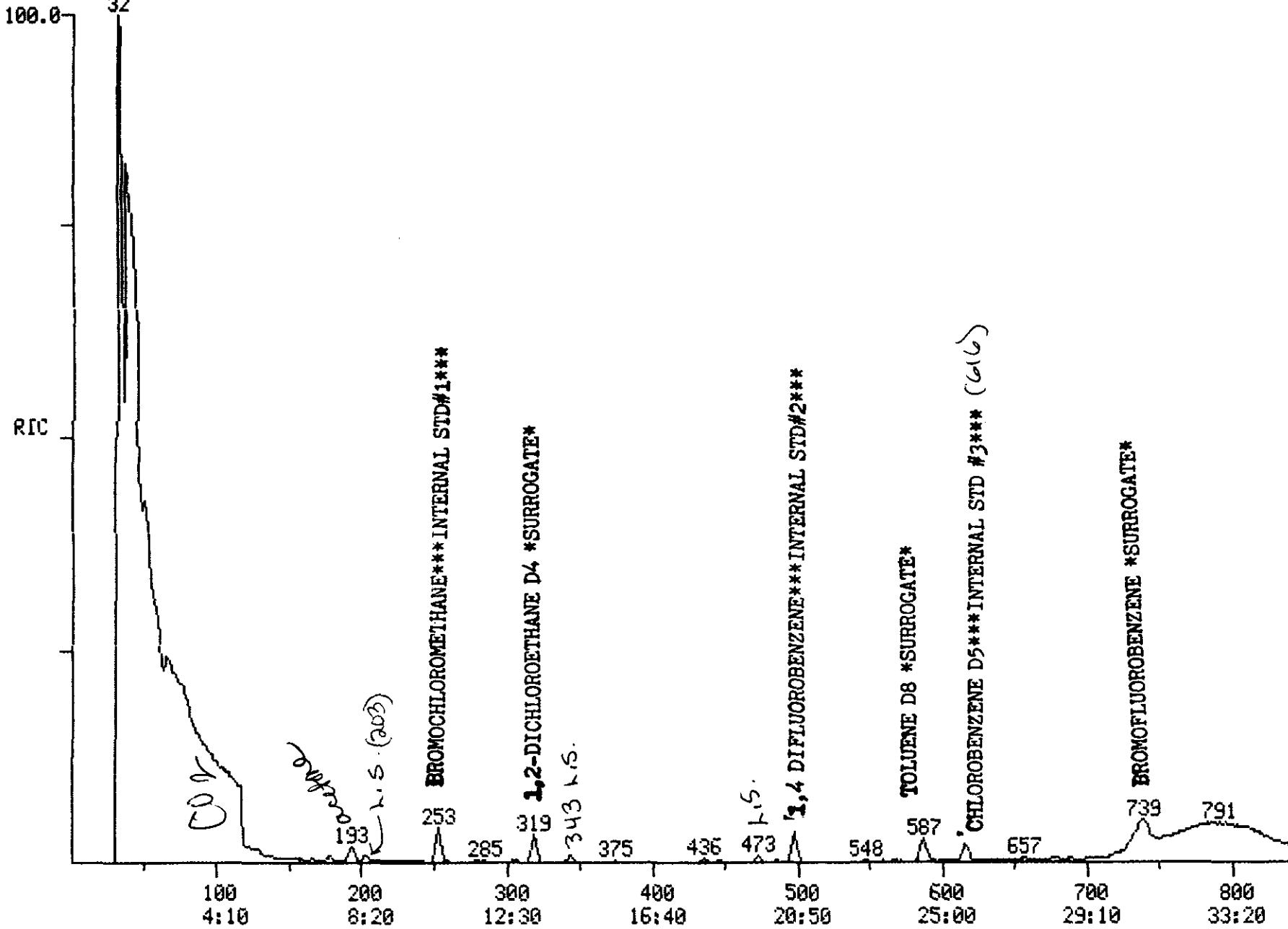
SAMPLE: CLP, UVERS CDM, 2536, 1, L, S, 16426, U, , 420.1.0, B2, 5G/5ML,,

COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

RANGE: G 1, 840 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

32

765952.
765952.17

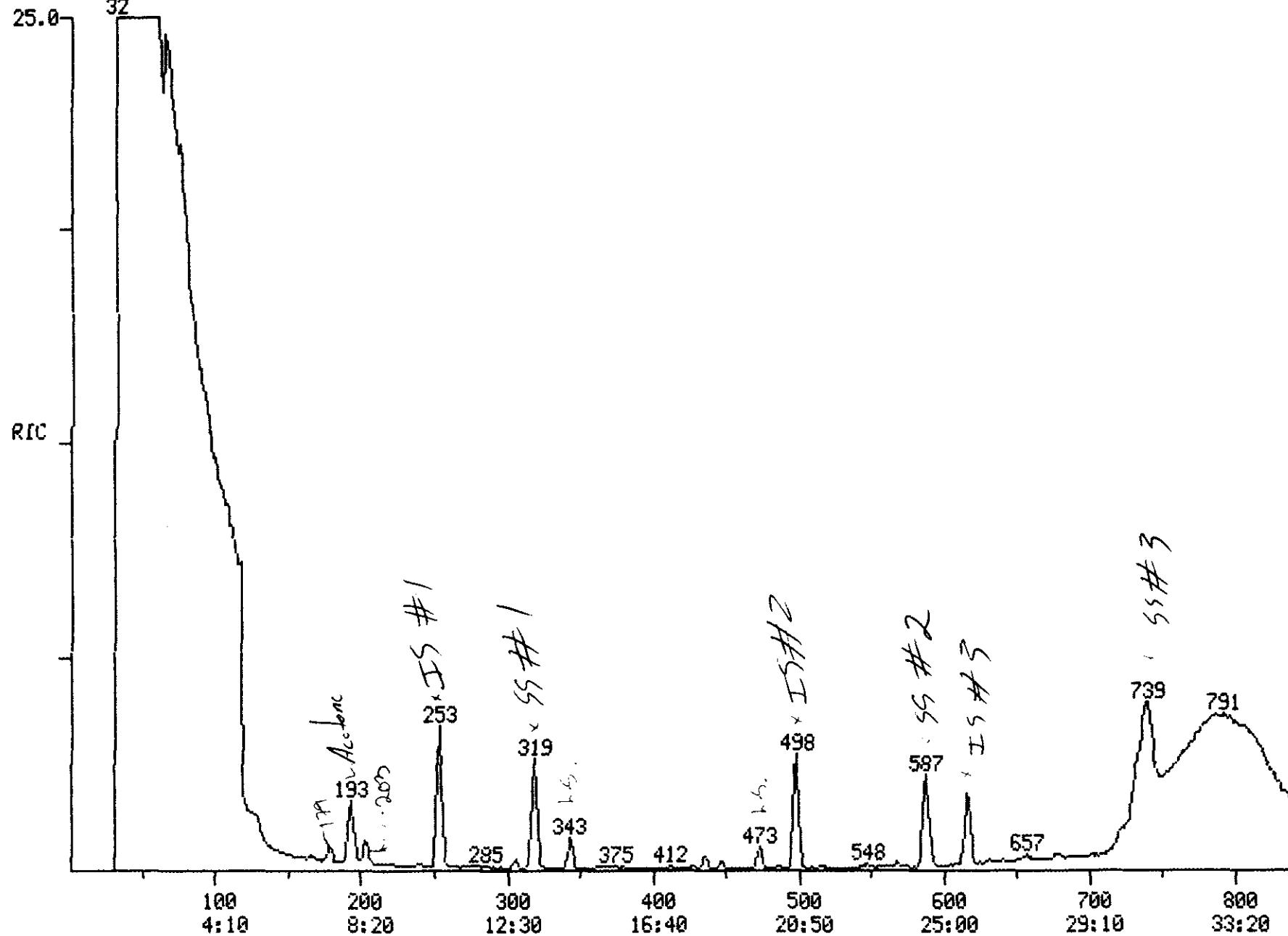


ORIGINAL

RIC
04/27/90 15:14:00
SAMPLE: CLP, VERSCDM, 2536, 1, L, S, 16426, U, , 420.1.0, 82, 5G/5ML,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
RANGE: G 1, 840 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: U3420 #1
CALI: U3420 #2
SCANS 1 TO 840

191488.
1000018



SCAN
TIME

Quantitation Report File: U3420

Data: U3420.TI

04/27/90 16:14:00

Sample: CLP, VERSCDM, 2536, 1, L, S, 16426, V, , 420, 1.0, B2, 5G/5ML.,

Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

Formula: - Instrument: U Weight: 0.011

Submitted by: VERSAR Analyst: MAT Acct. No.: 420.1.0

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)

Resp. fac. from Library Entry

30.0 %

No	Name
1	C101 BROMOCHLOROMETHANE **IS#1**
2	C010 CHLOROMETHANE
3	C015 BROMOMETHANE
4	C020 VINYL CHLORIDE
5	C025 CHLOROETHANE
6	C030 METHYLENE CHLORIDE
7	C035 ACETONE
8	C040 CARBON DISULFIDE
9	C045 1,1-DICHLOROETHENE
10	C050 1,1-DICHLOROETHANE
11	C053 1,2-DICHLOROETHENE (TOTAL)
12	C060 CHLOROFORM
13	C065 1,2-DICHLOROETHANE
14	CS15 1,2-DICHLOROETHANE-D4 **SS#1**
15	C043 TRICHLOROFUOROMETHANE
16	CI10 1,4-DIFLUOROBENZENE **IS#2**
17	C110 2-BUTANONE
18	C125 VINYL ACETATE
19	C120 CARBON TETRACHLORIDE
20	C130 BROMODICHLOROMETHANE
21	C140 1,2-DICLOROPROPANE
22	C145 CIS-1,3-DICLOROPROPENE
23	C150 TRICHLOROETHENE
24	C165 BENZENE
25	C155 DIBROMOCHLOROMETHANE
26	C160 1,1,2-TRICHLOROETHANE
27	C170 TRANS-1,3-DICLOROPROPENE
28	C180 BROMOFORM
29	C115 1,1,1-TRICHLOROETHANE
30	CS05 TOLUENE-D8 **SS#2**
31	CI20 CHLOROBENZENE-D5 **IS#3**
32	C205 4-METHYL-2-PENTANONE
33	C210 2-HEXANONE
34	C220 TETRACHLOROETHENE
35	C225 1,1,2,2-TETRACHLOROETHANE
36	C230 TOLUENE
37	C235 CHLOROBENZENE
38	C240 ETHYLBENZENE
39	C245 STYRENE
40	C250 ORTHO & PARA XYLENE
41	C251 META-XYLENE
42	CS10 BROMOFLUOROBENZENE **SS#3**

all (Same as low
2nd SS rec. high on

ready

100019

u34201

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	128	253	10:32	1	1.000	A BB	23936.	50.000 UG/L*	12.10
2	NOT FOUND						(low)		
3	NOT FOUND								
4	NOT FOUND								
5	NOT FOUND								
6	84	179	7:27	1	0.708	A BB	2411.	9.971 UG/L	0.96
7	43	193	8:02	1	0.763	A BV	25863.	109.293 UG/L	26.44
8	76	217	9:02	1	0.858	A BB	172.	0.458 UG/L	0.11
9	NOT FOUND							BRL	
10	NOT FOUND								
11	NOT FOUND								
12	NOT FOUND								
13	NOT FOUND								
14	65	319	13:17	1	1.261	A BB	34820.	36.217 UG/L%	8.76
15	NOT FOUND								
16	114	498	20:45	16	1.000	A BB	50685.	50.000 UG/L*	12.10
17	72	319	13:17	1	1.261	A BB	600.	6.355 UG/L	1.54
18	NOT FOUND							BRL	
19	NOT FOUND								
20	NOT FOUND								
21	NOT FOUND								
22	NOT FOUND								
23	NOT FOUND								
24	NOT FOUND								
25	NOT FOUND								
26	NOT FOUND								
27	NOT FOUND								
28	NOT FOUND								
29	NOT FOUND								
30	98	587	24:27	31	0.953	A BB	36239.	62.515 UG/L%	15.12
31	117	616	25:40	31	1.000	A BB	25216.	50.000 UG/L*	12.10
32	43	518	21:35	31	0.841	A BB	464.	1.269 UG/L	0.31
33	NOT FOUND							BRL	
34	NOT FOUND								
35	NOT FOUND								
36	92	591	24:37	31	0.959	A BB	320.	0.995 UG/L	0.24
37	NOT FOUND							BRL	
38	NOT FOUND								
39	NOT FOUND								
40	NOT FOUND								
41	NOT FOUND								
42	95	732	30:30	31	1.188	A BB	16081.	42.333 UG/L%	10.24

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
1	10:30	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
2	2:10		0.206						
3	3:20		0.317						
4	4:12		0.401						
5	5:15		0.500						
6	7:25	1.01	0.706	1.00	3.97	50.00	0.101	1.268	0.08
7	8:00	1.01	0.762	1.00	109.29	50.00	1.081	0.494	2.19
8	9:00	1.00	0.857	1.00	0.46	50.00	0.007	0.784	0.01
9	10:05		0.960						
10	11:22		1.083						
11	12:02		1.147						
12	12:37		1.202						

100020

U3420

(11) (10)

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
13	13: 20		1. 270						
14	13: 15	1. 00	1. 262	1. 00	36. 22	50. 00	1. 455	2. 008	0. 72
15	9: 27		0. 901						
16	20: 42	1. 00	1. 000	1. 00	50. 00	50. 00	1. 000	1. 000	1. 00
17	13: 15	1. 00	1. 262	1. 00	6. 36	50. 00	0. 025	0. 197	0. 13
18	15: 07		0. 730						
19	15: 05		0. 728						
20	15: 30		0. 748						
21	16: 55		0. 817						
22	17: 07		0. 827						
23	17: 40		0. 853						
24	18: 12		0. 879						
25	18: 17		0. 883						
26	18: 25		0. 889						
27	18: 22		0. 887						
28	20: 57		1. 012						
29	14: 42		0. 710						
30	24: 27	1. 00	0. 953	1. 00	62. 51	50. 00	1. 437	1. 149	1. 25
31	25: 40	1. 00	1. 000	1. 00	50. 00	50. 00	1. 000	1. 000	1. 00
32	21: 30	1. 00	0. 838	1. 00	1. 27	50. 00	0. 018	0. 725	0. 03
33	23: 02		0. 898						
34	23: 17		0. 907						
35	23: 17		0. 907						
36	24: 37	1. 00	0. 959	1. 00	0. 99	50. 00	0. 013	0. 638	0. 02
37	25: 47		1. 005						
38	27: 57		1. 089						
39	32: 25		1. 263						
40	33: 47		1. 317						
41	32: 42		1. 274						
42	30: 45	0. 99	1. 198	0. 99	42. 33	50. 00	0. 638	0. 753	0. 85

100021

MASS SPECTRUM

04/27/90 16:14:00 + 8:02

SAMPLE: CLP,VERSCDM,2536,1,L,S,16426,U,,420.1.0,B2,5G/5ML,,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
** NAME: C035 ACETONE

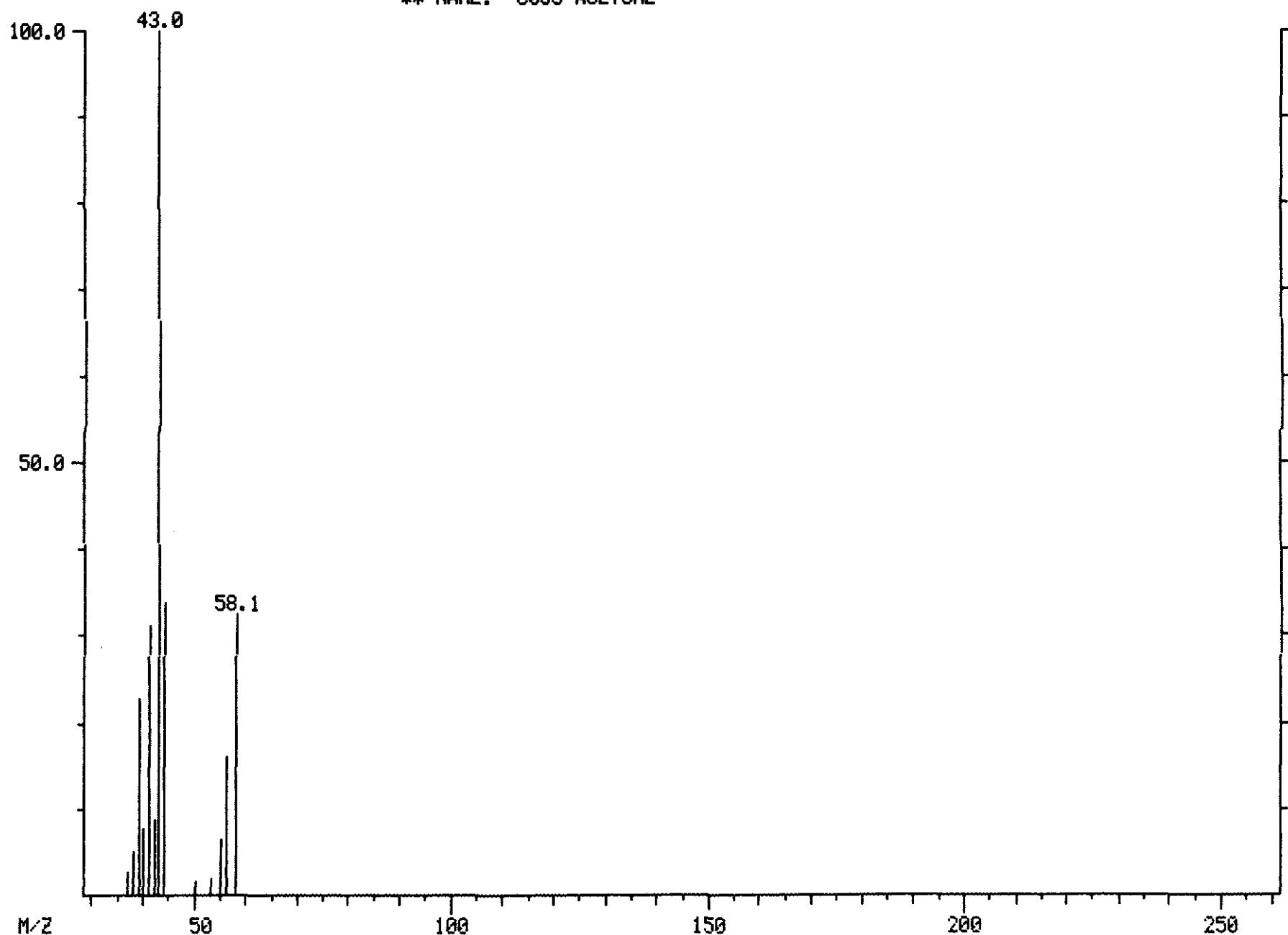
DATA: U3420 #193

CALI: U3420 #2

BASE M/Z: 43

RIC: 15680.

000222
5776.



MASS SPECTRUM

04/27/90 16:14:00 + 8:02

SAMPLE: CLP, VERSCDM, 2536, 1, L, S, 16426, U, , 420.1.0, B2, 5G/5ML,,

COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

** NAME: C035 ACETONE

ENHANCED (S 15B 2N 0T)

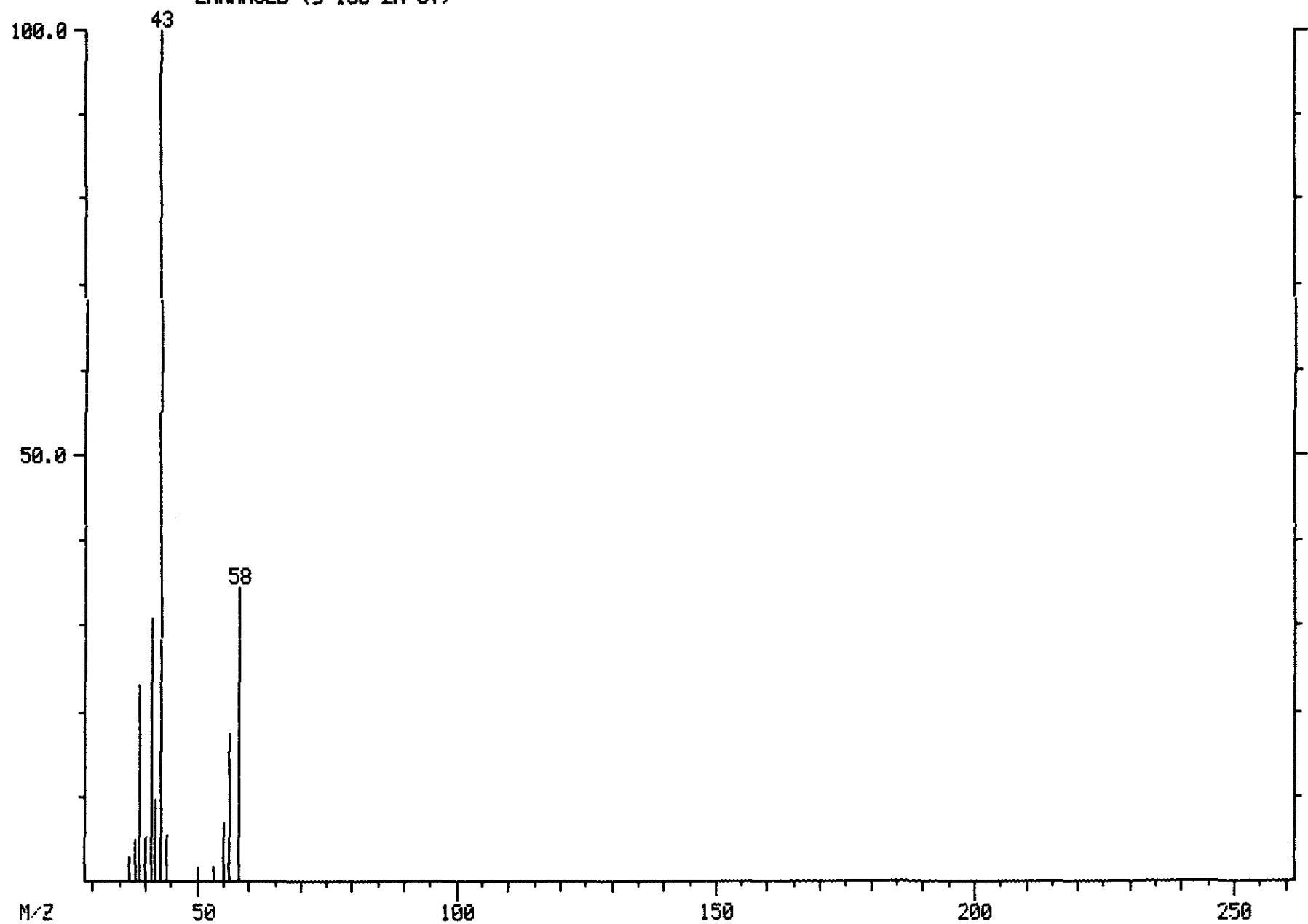
DATA: U3420 #193

CALI: U3420 #2

BASE M/Z: 43

RIC: 13168.

100023



MASS SPECTRUM
04/27/90 8:21:00 + 8:00

SAMPLE: CLP,,,VSTD 50,L,S,22554,V,CC-050,,5ML,,

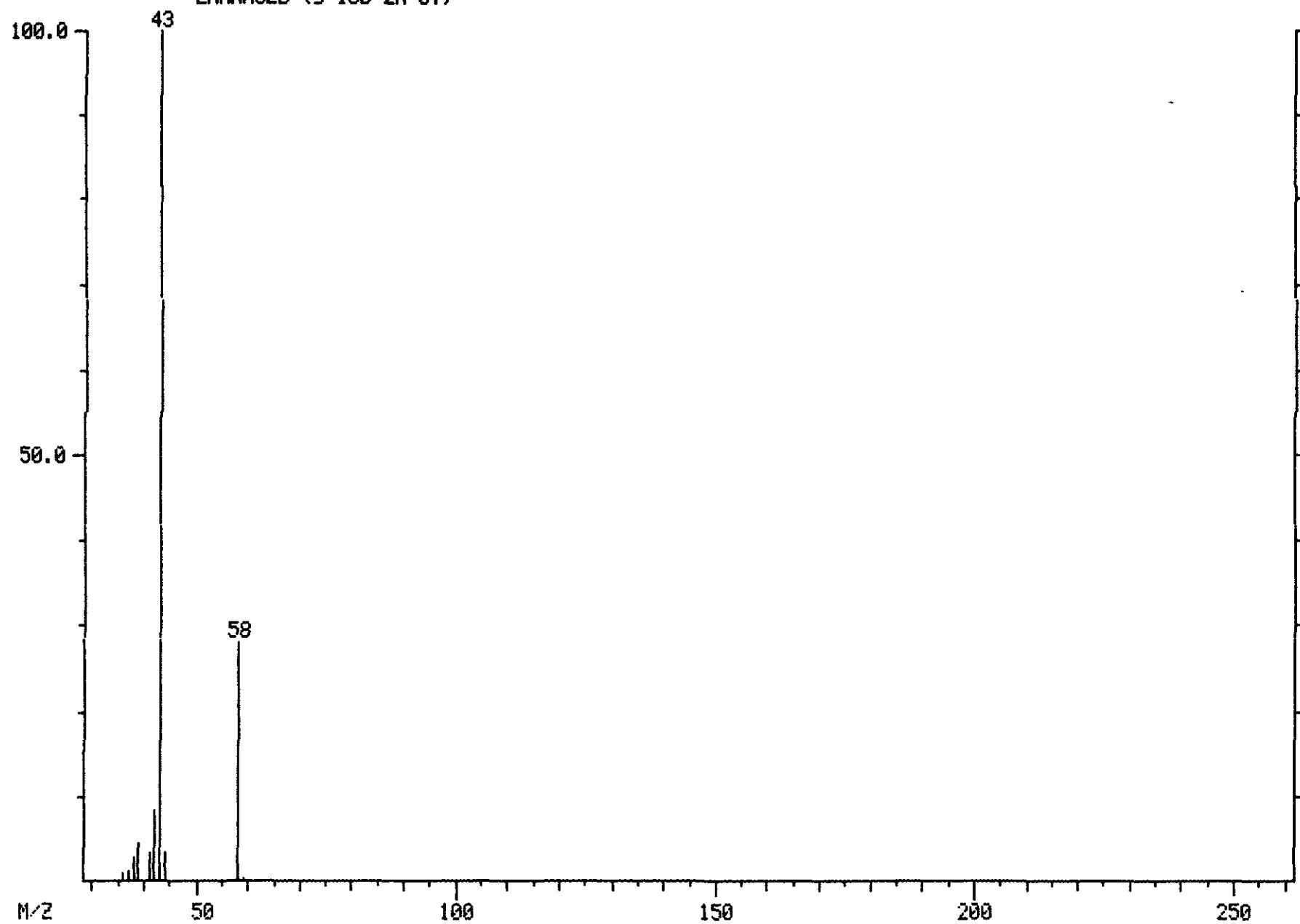
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

** NAME: C035 ACETONE

ENHANCED (S 15B 2N 0T)

DATA: U3411 #192
CALI: U3411 #2

BASE M/Z: 43
RIC: 12368.



100024

Library Search Data: U3420 # 203 Base m/z: 43
 04/27/90 16:14:00 + 8:27 Cali: U3420 # 2 RIC: 4127~
 Sample: CLP, VERSCDM, 2536, 1, L, S, 16426, V., 420. 1. O, B2, 5G/5ML,
 Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
 Enhanced (S 15B 2N OT)

42222 spectra in LIBRARYNB searched for maximum PURITY
 78 matched at least 5 of the 6 largest peaks in the unknown

Rank In. Name

1	94 BUTANE
2	93 PROPANE, 2-METHYL-
3	1261 1-PROPEN-2-OL, ACETATE
4	847 PROPANE, 1-CHLORO-2-METHYL-
5	1204 PROPANE, 1-ISOCYANATO-2-METHYL-

Rank	Formula	M. Wt	B. Pk	Purity	Fit	RFit
1	C4. H10	58	43	891	911	965~
2	C4. H10	58	43	837	856	901
3	C5. H8. O2	100	43	756	773	916
4	C4. H9. CL	92	43	731	850	833
5	C5. H9. O. N	99	43	712	712	866

Rank	Ret. Time	B. P. Int.	US. Par. 1	US. Par. 2	C. A. S. #
1	—	—	—	—	106-97-8~
2	—	—	—	—	75-28-5
3	—	—	—	—	108-22-5
4	—	—	—	—	513-36-0
5	—	—	—	—	1873-29-6

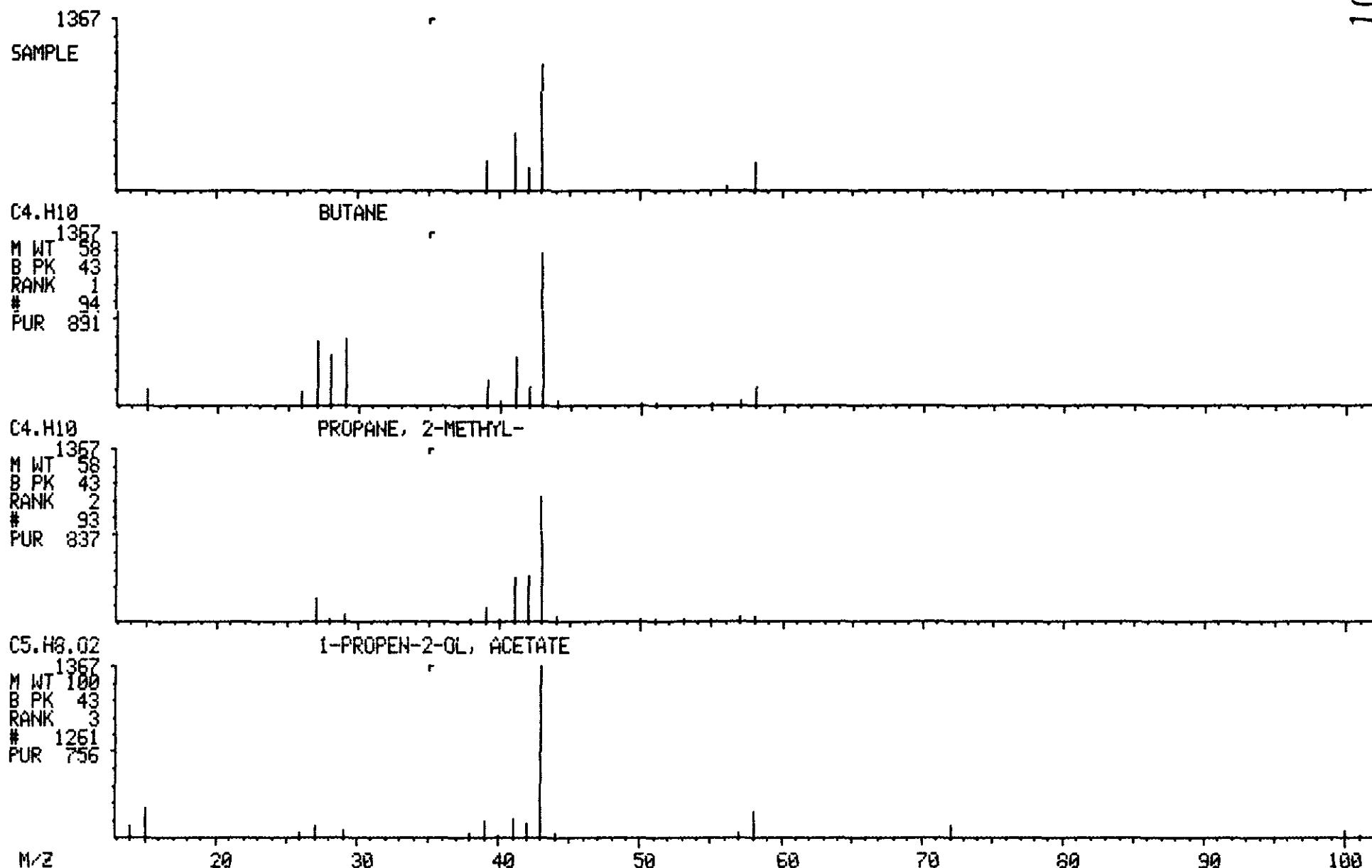
~~4127~~
 30365 x 50 = 6.79

10025

LIBRARY SEARCH
04/27/90 15:14:00 + 8:27
SAMPLE: CLP,VERSCDM,2536,1,L,S,16426,U,,420.1.0,B2,5G/5ML,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
ENHANCED (S 158 2N 0T)

DATA: U3420 # 203
CALI: U3420 # 2

BASE M/Z: 43
RIC: 4127.



Library Search Data: U3420 # 343 Base m/z: 43
 04/27/90 16:14:00 + 14:17 Cali: U3420 # 2 RIC: 6455✓
 Sample: CLP, VERSCDM, 2536, 1, L, S, 16426, V, , 420. 1. 0, B2, 5G/5ML.,
 Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
 Enhanced (S 15B 2N OT)

42222 spectra in LIBRARYNB searched for maximum PURITY
 58 matched at least 6 of the 9 largest peaks in the unknown

Rank In. Name

- 1 265 PENTANE (ACN)(DOT)
- 2 253 OXIRANE, 2,2-DIMETHYL-
- 3 93 PROPANE, 2-METHYL-
- 4 249 1-PROPENE, 2-METHOXY-
- 5 2512 ETHER, 1-BUTYLVINYL METHYL

Rank	Formula	M. Wt	B. Pk	Purity	Fit	RFit
1	C5. H12	72	43	974	974	991 ✓
2	C4. H8. O	72	41	905	914	935
3	C4. H10	58	43	827	919	865
4	C4. H8. O	72	42	801	821	833
5	C7. H14. O	114	42	783	794	871

Rank	Ret. Time	B. P. Int.	US. Par. 1	US. Par. 2	C. A. S. #
1	—	—	—	—	109-66-0 ✓
2	—	—	—	—	558-30-5
3	—	—	—	—	75-28-5
4	—	—	—	—	116-11-0
5	—	—	—	—	16519-66-7

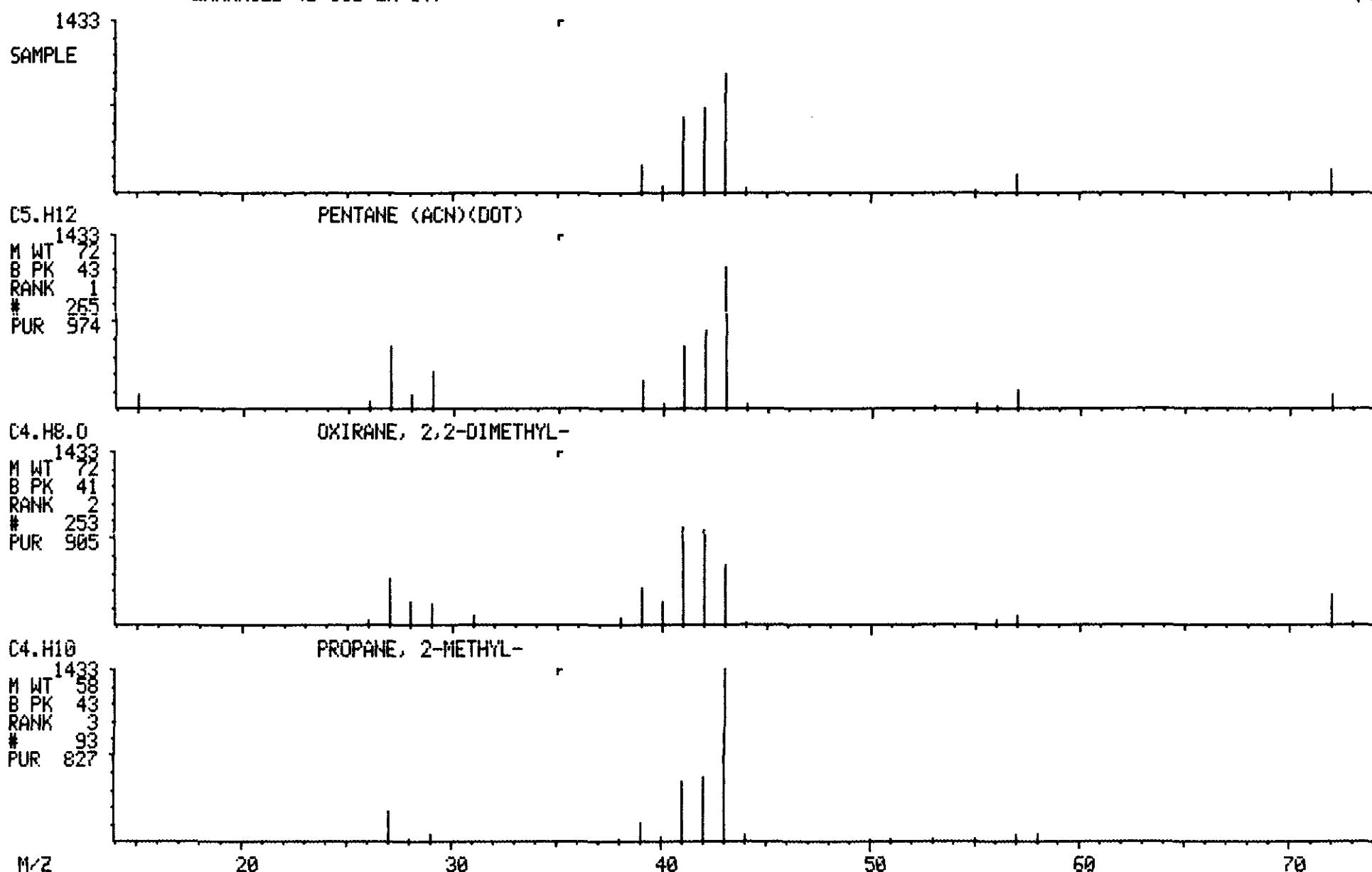
~~6455~~
~~30368~~ 10.62

100027

LIBRARY SEARCH
04/27/90 16:14:00 + 14:17
SAMPLE: CLP, VERSCDM, 2536, 1, L, S, 16426, U,, 420.1.0, B2, 5G/5ML,,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
ENHANCED (S 15B 2N 0T)

DATA: U3420 # 343
CALI: U3420 # 2

BASE M/Z: 43
RIC: 6455.



1000028

Library Search Data: U3420 # 473 Base m/z: 57
 04/27/90 16:14:00 + 19:42 Cali: U3420 # 2 RIC: 4559 ✓
 Sample: CLP, VERSCDM, 2536, 1, L, S, 16426, V, , 420. 1. 0, B2, 5G/5ML.,
 Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN RI
 Enhanced (S 15B 2N OT)

42222 spectra in LIBRARYNB searched for maximum PURITY
 97 matched at least 6 of the 8 largest peaks in the unknown

Rank In. Name

- 1 664 HEXANE (DOT)
- 2 662 PENTANE, 3-METHYL-
- 3 633 PROPANAL, 2, 2-DIMETHYL-
- 4 1548 PROPANE, 2-METHYL-1-NITRO-
- 5 1370 PENTANE, 2, 4-DIMETHYL-

Rank	Formula	M. Wt	B. Pk	Purity	Fit	RFit
1	C6. H14	86	57	924	924	987 ✓
2	C6. H14	86	57	857	857	980
3	C5. H10. O	86	57	797	797	867
4	C4. H9. O2. N	103	41	782	882	802
5	C7. H16	100	43	768	768	888

Rank	Ret. Time	B. P. Int.	US. Par. 1	US. Par. 2	C. A. S. #
1	—	—	—	—	110-54-3 ✓
2	—	—	—	—	96-14-0
3	—	—	—	—	630-19-3
4	—	—	—	—	625-74-1
5	—	—	—	—	108-08-7

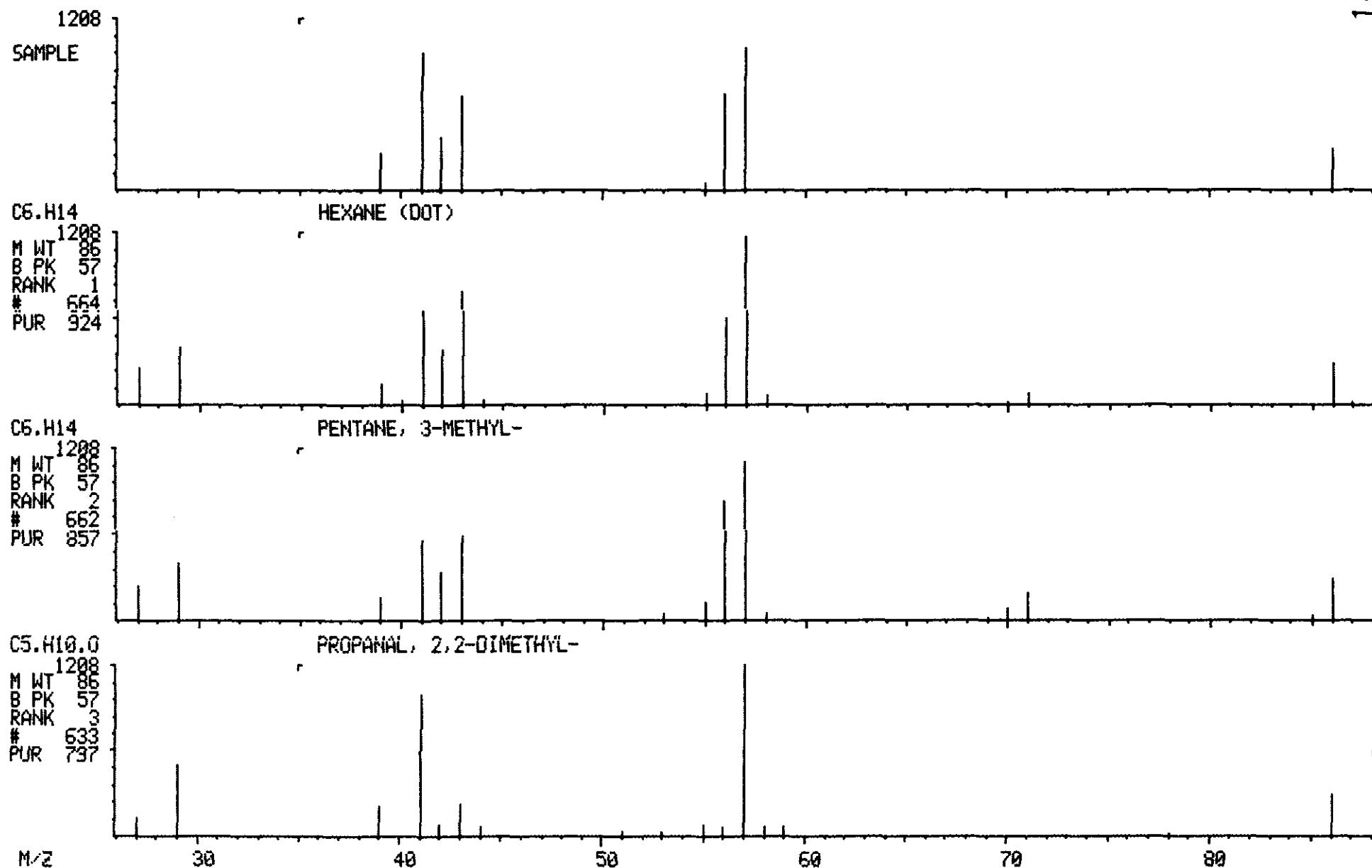
$\frac{4559}{25216} \times 50 = 9.03$

100029

LIBRARY SEARCH
04/27/90 16:14:00 + 19:42
SAMPLE: CLP,VERSCDM,2536,1,L,S,16426,U,,420.1.0,B2,5G/5ML,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
ENHANCED (S 15B 2N 0T)

DATA: U3420 # 473
CALI: U3420 # 2

BASE M/Z: 57
RIC: 4559.



100030

Library Search Data: U3420 # 791 Base m/z: 128
 04/27/90 16:14:00 + 32:57 Cali: U3420 # 2 RIC: 17471.
 Sample: CLP, VERSCDM, 2536, 1, L, S, 16426, V, , 420, 1, 0, B2, 5G/5ML.,,
 Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
 # 791 - # 838 - # 750 X1.00

42222 spectra in LIBRARYNB searched for maximum PURITY
 123 matched at least 5 of the 16 largest peaks in the unknown

PAH

Rank In. Name

- 1 3887 AZULENE
- 2 3886 NAPHTHALENE (ACN)(DOT)
- 3 3888 1H-INDENE, 1-METHYLENE-
- 4 8736 BICYCLO[4.4.1]UNDECA-1,3,5,7,9-PENTAEN-11-ONE
- 5 4116 1,4-BENZENEDICARBONITRILE

Rank	Formula	M. Wt	B. Pk	Purity	Fit	RFit
1	C10. H8	128	128	946	997	946
2	C10. H8	128	128	941	997	941
3	C10. H8	128	128	937	996	937
4	C11. H8. O	156	128	791	887	882
5	C8. H4. N2	128	128	729	895	749

Rank	Ret. Time	B. P. Int.	US. Par. 1	US. Par. 2	C. A. S. #
1	---	---	---	---	275-51-4
2	---	---	---	---	91-20-3
3	---	---	---	---	2471-84-3
4	---	---	---	---	36628-80-5
5	---	---	---	---	623-26-7

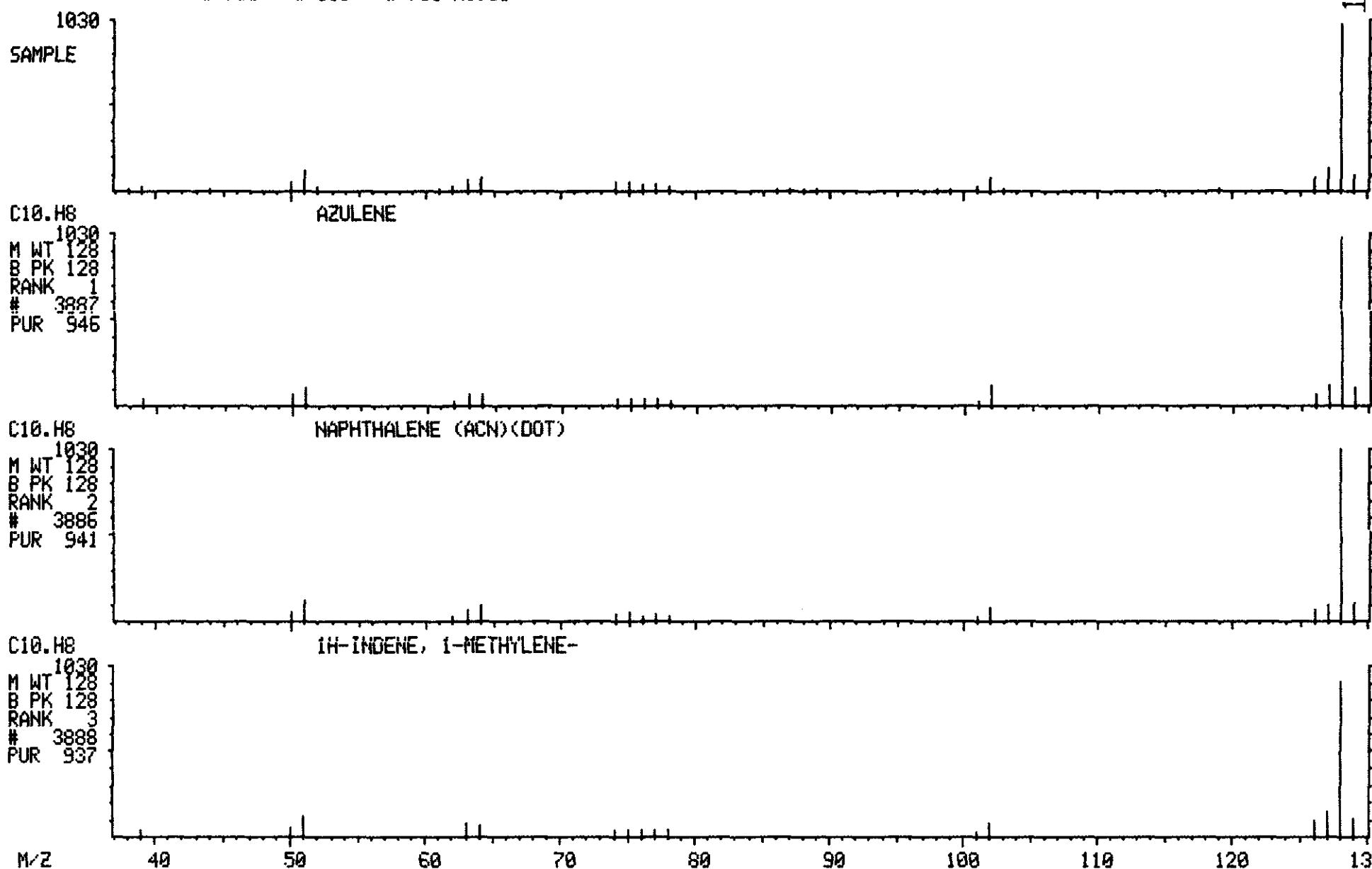
$$\frac{17471}{15568} \times 50 = 56.11$$

100031

LIBRARY SEARCH
04/27/90 16:14:00 + 32:57
SAMPLE: CLP, VERSCDM, 2536, 1, L, 5, 16426, U,, 420.1.0, B2, 5G/5ML,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
791 - # 838 - # 750 X1.00

DATA: U3420 # 791
CALI: U3420 # 2

BASE M/Z: 128
RIC: 17471.



100032

100033

ANALYST: CHECK BASE M/Z AND RIC AMOUNT TO INSURE NO CONTAMINATION

*****INTERNAL STANDARD RIC REPORT<-->
Mass L15t Data: U3420 # 253 Base m/z: 130
04/27/90 16:14:00 + 10:32 Cai1: U3420 # 2 RIC: 30368.
Sample: CLP, VERSCDM, 2536, 1,L,S, 16426, V,, 420, 1,O,B2, 5G/5ML,
Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ BC/MIN
Enhanced (S 15B 2N OT)
*****INTERNAL STANDARD RIC#1*****
Mass L15t Data: U3420 # 253 Base m/z: 130
04/27/90 16:14:00 + 10:32 Cai1: U3420 # 2 RIC: 30368.
Sample: CLP, VERSCDM, 2536, 1,L,S, 16426, V,, 420, 1,O,B2, 5G/5ML,
Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ BC/MIN
Enhanced (S 15B 2N OT)
*****INTERNAL STANDARD RIC#2*****
Mass L15t Data: U3420 # 498 Base m/z: 114
04/27/90 16:14:00 + 20:45 Cai1: U3420 # 2 RIC: 25216.
Sample: CLP, VERSCDM, 2536, 1,L,S, 16426, V,, 420, 1,O,B2, 5G/5ML,
Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ BC/MIN
Enhanced (S 15B 2N OT)
*****INTERNAL STANDARD RIC#3*****
Mass L15t Data: U3420 # 616 Base m/z: 117
04/27/90 16:14:00 + 25:40 Cai1: U3420 # 2 RIC: 15568.
Sample: CLP, VERSCDM, 2536, 1,L,S, 16426, V,, 420, 1,O,B2, 5G/5ML,
Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ BC/MIN
Enhanced (S 15B 2N OT)

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: VERSAR INC.

Contract: _____

2

Lab Code: VERSAR Case No.: R3-7

SAS No.: _____ SDG No.: 1

Matrix: (soil/water) SOIL

Lab Sample ID: 16427

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: U3393

Level: (low/med) LOW

Date Received: 04/19/90

* Moisture: not dec. 27

Date Analyzed: 04/26/90

Column: (pack/cap) PACK

Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG	Q
74-87-3-----	Chloromethane	14	U
74-83-9-----	Bromomethane	14	U
75-01-4-----	Vinyl chloride	14	U
75-00-3-----	Chloroethane	14	U
75-09-2-----	Methylene chloride	7	U
67-64-1-----	Acetone	14	U
75-15-0-----	Carbon disulfide	7	U
75-35-4-----	1,1-Dichloroethene	7	U
75-34-3-----	1,1-Dichloroethane	7	U
540-59-0-----	1,2-Dichloroethene (total)	7	U
67-66-3-----	Chloroform	7	U
107-06-2-----	1,2-Dichloroethane	7	U
78-93-3-----	2-Butanone	14	U
71-55-6-----	1,1,1-Trichloroethane	7	U
56-23-5-----	Carbon tetrachloride	7	U
108-05-4-----	Vinyl acetate	14	U
75-27-4-----	Bromodichloromethane	7	U
78-87-5-----	1,2-Dichloropropane	7	U
10061-01-5-----	cis-1,3-Dichloropropene	7	U
79-01-6-----	Trichloroethene	7	U
124-48-1-----	Dibromochloromethane	7	U
79-00-5-----	1,1,2-Trichloroethane	7	U
71-43-2-----	Benzene	7	U
10061-02-6-----	Trans-1,3-dichloropropene	7	U
75-25-2-----	Bromoform	7	U
108-10-1-----	4-Methyl-2-pentanone	14	U
591-78-6-----	2-Hexanone	14	U
127-18-4-----	Tetrachloroethene	7	U
79-34-5-----	1,1,2,2-Tetrachloroethane	7	U
108-88-3-----	Toluene	7	U
108-90-7-----	Chlorobenzene	7	U
100-41-4-----	Ethylbenzene	7	U
100-42-5-----	Styrene	7	U
1330-20-7-----	Total xylenes	7	U

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: VERSAR INC.

Contract: _____

2

Lab Code: VERSAR Case No.: R3-7

SAS No.: _____ SDG No.: 1 _____

Matrix: (soil/water) SOIL

Lab Sample ID: 16427 _____

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: U3393 _____

Level: (low/med) LOW

Date Received: 04/19/90

* Moisture: not dec. 27

Date Analyzed: 04/26/90

Column (pack/cap) PACK

Dilution Factor: 1.0 _____

N^o ber TICs found: 0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

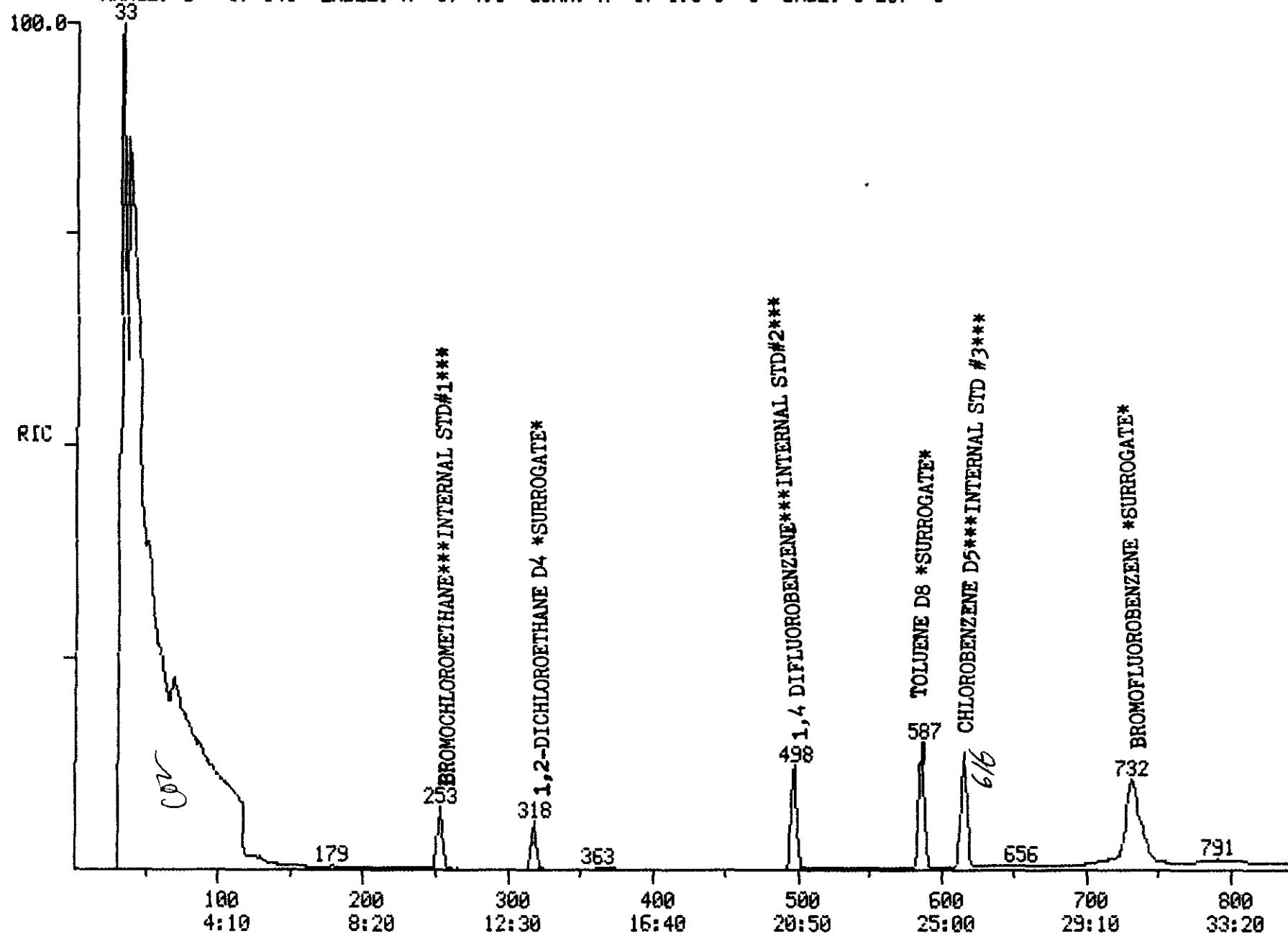
RIC
04/26/90 14:31:00

DATA: U3393 #1
CALI: U3393 #2

SCANS 1 TO 840

SAMPLE: CLP, VERSCDM, 2536, B2, L, S, 16427, U, , 420.1.0, B2, 5G/SML.,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

RANGE: G 1, 840 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3



732160.
00036

Quantitation Report File: U3393

Data: U3393.TI

04/26/90 14:31:00

Sample: CLP, VERSCDM, 2536, B2, L, S, 16427, V., 420.1.0, B2, 5G/5ML, ,

Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ BC/MIN

Formula: -

Instrument: U

Weight: 0.011

Submitted by: VERSAR

Analyst: MAT

Acct. No.: 420.1.0

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)

Resp. fac. from Library Entry

INT STD + SURROG. Comp.

No	Name
1	C101 BROMOCHLOROMETHANE **IS#1**
2	C010 CHLOROMETHANE
3	C015 BROMOMETHANE
4	C020 VINYL CHLORIDE
5	C025 CHLOROETHANE
6	C030 METHYLENE CHLORIDE
7	C035 ACETONE
8	C040 CARBON DISULFIDE
9	C045 1,1-DICHLOROETHENE
10	C050 1,1-DICHLOROETHANE
11	C053 1,2-DICHLOROETHENE (TOTAL)
12	C060 CHLOROFORM
13	C065 1,2-DICHLOROETHANE
14	CS15 1,2-DICHLOROETHANE-D4 **SS#1**
15	C043 TRICHLOROFLUOROMETHANE
16	CI10 1,4-DIFLUOROBENZENE **IS#2**
17	C110 2-BUTANONE
18	C125 VINYL ACETATE
19	C120 CARBON TETRACHLORIDE
20	C130 BROMODICHLOROMETHANE
21	C140 1,2-DICHLOROPROPANE
22	C145 CIS-1,3-DICHLOROPROPENE
23	C150 TRICHLOROETHENE
24	C165 BENZENE
25	C155 DIBROMOCHLOROMETHANE
26	C160 1,1,2-TRICHLOROETHANE
27	C170 TRANS-1,3-DICHLOROPROPENE
28	C180 BROMOFORM
29	C115 1,1,1-TRICHLOROETHANE
30	CS05 TOLUENE-D8 **SS#2**
31	CI20 CHLOROBENZENE-D5 **IS#3**
32	C205 4-METHYL-2-PENTANONE
33	C210 2-HEXANONE
34	C220 TETRACHLOROETHENE
35	C225 1,1,2,2-TETRACHLOROETHANE
36	C230 TOLUENE
37	C235 CHLOROBENZENE
38	C240 ETHYLBENZENE
39	C245 STYRENE
40	C250 ORTHO & PARA XYLENE
41	C251 META-XYLENE
42	CS10 BROMOFLUOROBENZENE **SS#3**

M.A. Templeton
4/26/90
27.5 %

Ready

100037

U3393

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	128	253	10:32	1	1.000	A BB	38573.	50.000 UG/L*	17.56
2	NOT FOUND						OK.		
3	NOT FOUND						✓		
4	NOT FOUND								
5	NOT FOUND								
6	84	180	7:30	1	0.711	A BB	1258.	+ 189 UG/L	0.42
7	43	193	8:02	1	0.763	A BB	179.	- 0.319 UG/L	0.11
8	NOT FOUND								
9	NOT FOUND								
10	NOT FOUND								
11	NOT FOUND								
12	NOT FOUND								
13	NOT FOUND								
14	65	318	13:15	1	1.257	A BB	61885.	35.065 UG/L%	12.31 70%
15	NOT FOUND								
16	114	498	20:45	16	1.000	A BB	165489.	✓ 50.000 UG/L*	17.56
17	NOT FOUND								
18	NOT FOUND								
19	NOT FOUND								
20	NOT FOUND								
21	NOT FOUND								
22	NOT FOUND								
23	NOT FOUND								
24	NOT FOUND								
25	NOT FOUND								
26	NOT FOUND								
27	NOT FOUND								
28	NOT FOUND								
29	NOT FOUND								
30	98	586	24:25	31	0.951	A BB	186603.	✓ 53.162 UG/L%	18.67 106%
31	117	616	25:40	31	1.000	A BB	145981.	✓ 50.000 UG/L*	17.56
32	NOT FOUND								
33	43	556	23:10	31	0.903	A VV	1804.	- 0.814 UG/L	0.29
34	NOT FOUND								
35	NOT FOUND								
36	92	591	24:37	31	0.959	A BB	1121.	- 0.559 UG/L	0.20
37	NOT FOUND								
38	NOT FOUND								
39	NOT FOUND								
40	106	796	33:10	31	1.292	A BB	300.	- 0.215 UG/L	0.08
41	NOT FOUND								
42	95	731	30:27	31	1.187	A BB	110417.	43.521 UG/L%	15.28 87

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
1	10:32	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
2	2:12		0.209						
3	3:22		0.320						
4	4:15		0.403						
5	5:15		0.498						
6	7:27	1.01	0.708	1.01	1.19	50.00	0.033	1.372	0.02
7	8:02	1.00	0.763	1.00	0.32	50.00	0.005	0.727	0.01
8	9:02		0.858						
9	10:07		0.960						
10	11:25		1.083						
11	12:05		1.146						
12	12:40		1.202						

100038

W3393

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R.Fac	R.Fac(L)	Ratio
13	13:22		1. 269						
14	13:17	1. 00	1. 261	1. 00	35. 06	50. 00	1. 604	2. 288	0. 70
15	9:30		0. 901						
16	20:45	1. 00	1. 000	1. 00	50. 00	50. 00	1. 000	1. 000	1. 00
17	13:17		1. 261						
18	15:10		0. 731						
19	15:07		0. 729						
20	15:32		0. 749						
21	16:57		0. 817						
22	17:10		0. 827						
23	17:42		0. 853						
24	18:12		0. 878						
25	18:20		0. 884						
26	18:25		0. 888						
27	18:25		0. 888						
28	21:00		1. 012						
29	14:42		0. 709						
30	24:27	1. 00	0. 953	1. 00	53. 16	50. 00	1. 278	1. 202	1. 06
31	25:40	1. 00	1. 000	1. 00	50. 00	50. 00	1. 000	1. 000	1. 00
32	21:30		0. 838						
33	23:02	1. 01	0. 898	1. 01	0. 81	50. 00	0. 012	0. 759	0. 02
34	23:17		0. 907						
35	23:17		0. 907						
36	24:37	1. 00	0. 959	1. 00	0. 56	50. 00	0. 008	0. 686	0. 01
37	25:47		1. 005						
38	27:50		1. 084						
39	32:00		1. 247						
40	33:17	1. 00	1. 297	1. 00	0. 22	50. 00	0. 002	0. 477	0. 00
41	32:15		1. 256						
42	30:27	1. 00	1. 187	1. 00	43. 52	50. 00	0. 756	0. 869	0. 87

100039

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: VERSAR INC.

Contract: _____

3

Lab Code: VERSAR Case No.: R3-7

SAS No.: _____ SDG No.: 1

Matrix: (soil/water) SOIL

Lab Sample ID: 16428

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: U3394

Level: (low/med) LOW

Date Received: 04/19/90

% Moisture: not dec. 28

Date Analyzed: 04/26/90

Column: (pack/cap) PACK

Dilution Factor: 1.0

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

<u>74-87-3-----Chloromethane</u>	<u>14</u>	<u>U</u>
<u>74-83-9-----Bromomethane</u>	<u>14</u>	<u>U</u>
<u>75-01-4-----Vinyl chloride</u>	<u>14</u>	<u>U</u>
<u>75-00-3-----Chloroethane</u>	<u>14</u>	<u>U</u>
<u>75-09-2-----Methylene chloride</u>	<u>7</u>	<u>U</u>
<u>67-64-1-----Acetone</u>	<u>14</u>	<u>U</u>
<u>75-15-0-----Carbon disulfide</u>	<u>7</u>	<u>U</u>
<u>75-35-4-----1,1-Dichloroethene</u>	<u>7</u>	<u>U</u>
<u>75-34-3-----1,1-Dichloroethane</u>	<u>7</u>	<u>U</u>
<u>540-59-0-----1,2-Dichloroethene (total)</u>	<u>7</u>	<u>U</u>
<u>67-66-3-----Chloroform</u>	<u>7</u>	<u>U</u>
<u>107-06-2-----1,2-Dichloroethane</u>	<u>7</u>	<u>U</u>
<u>78-93-3-----2-Butanone</u>	<u>14</u>	<u>U</u>
<u>71-55-6-----1,1,1-Trichloroethane</u>	<u>7</u>	<u>U</u>
<u>56-23-5-----Carbon tetrachloride</u>	<u>7</u>	<u>U</u>
<u>108-05-4-----Vinyl acetate</u>	<u>14</u>	<u>U</u>
<u>75-27-4-----Bromodichloromethane</u>	<u>7</u>	<u>U</u>
<u>78-87-5-----1,2-Dichloropropane</u>	<u>7</u>	<u>U</u>
<u>10061-01-5-----cis-1,3-Dichloropropene</u>	<u>7</u>	<u>U</u>
<u>79-01-6-----Trichloroethene</u>	<u>7</u>	<u>U</u>
<u>124-48-1-----Dibromochloromethane</u>	<u>7</u>	<u>U</u>
<u>79-00-5-----1,1,2-Trichloroethane</u>	<u>7</u>	<u>U</u>
<u>71-43-2-----Benzene</u>	<u>7</u>	<u>U</u>
<u>10061-02-6-----Trans-1,3-dichloropropene</u>	<u>7</u>	<u>U</u>
<u>75-25-2-----Bromoform</u>	<u>7</u>	<u>U</u>
<u>108-10-1-----4-Methyl-2-pentanone</u>	<u>14</u>	<u>U</u>
<u>591-78-6-----2-Hexanone</u>	<u>14</u>	<u>U</u>
<u>127-18-4-----Tetrachloroethene</u>	<u>7</u>	<u>U</u>
<u>79-34-5-----1,1,2,2-Tetrachloroethane</u>	<u>7</u>	<u>U</u>
<u>108-88-3-----Toluene</u>	<u>7</u>	<u>U</u>
<u>108-90-7-----Chlorobenzene</u>	<u>7</u>	<u>U</u>
<u>100-41-4-----Ethylbenzene</u>	<u>7</u>	<u>U</u>
<u>100-42-5-----Styrene</u>	<u>7</u>	<u>U</u>
<u>1330-20-7-----Total xylenes</u>	<u>7</u>	<u>U</u>

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: VERSAR INC.

Contract: _____

3

Lab Code: VERSAR Case No.: R3-7

SAS No.: _____ SDG No.: 1 _____

Matrix: (soil/water) SOIL

Lab Sample ID: 16428 _____

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: U3394 _____

Level: (low/med) LOW

Date Received: 04/19/90

* Moisture: not dec. 28

Date Analyzed: 04/26/90

Column (pack/cap) PACK

Dilution Factor: 1.0 _____

Number TICs found: 0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

RIC
04/26/90 15:24:00

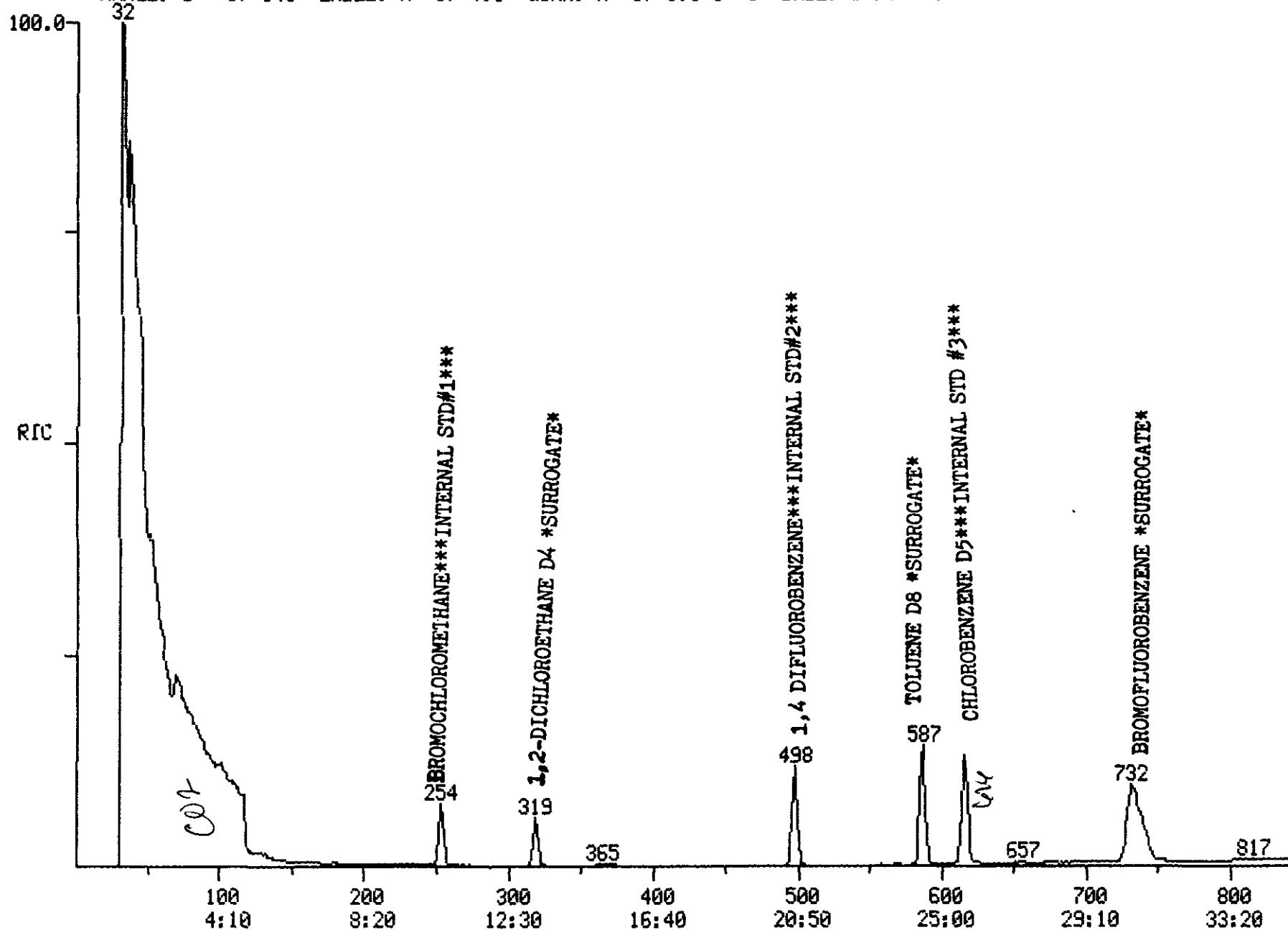
DATA: U3394 #1
CALI: U3394 #2

SCANS 1 TO 840

SAMPLE: CLP,VERSCM,2536,B3,L,S,16428,U,,420.1.0,B2,5G/5ML,,

COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

RANGE: G 1, 840 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3



Quantitation Report File: U3394

Data: U3394.TI

04/26/90 15:24:00

Sample: CLP, VERSCM, 2536, #3, L, S, 16428, V, , 420. 1. 0, B2, 5G/5ML.,

Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

Formula: - Instrument: U Weight: 0.011

Submitted by: VERSAR Analyst: MAT Acct. No.: 420. 1. 0

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)

Resp. fac. from Library Entry

28.4 %

No	Name
1	C101 BROMOCHLOROMETHANE **IS#1**
2	C010 CHLOROMETHANE
3	C015 BROMOMETHANE
4	C020 VINYL CHLORIDE
5	C025 CHLOROETHANE
6	C030 METHYLENE CHLORIDE
7	C035 ACETONE
8	C040 CARBON DISULFIDE
9	C045 1,1-DICHLOROETHENE
10	C050 1,1-DICHLOROETHANE
11	C053 1,2-DICHLOROETHENE (TOTAL)
12	C060 CHLOROFORM
13	C065 1,2-DICHLOROETHANE
14	CS15 1,2-DICHLOROETHANE-D4 **SS#1**
15	C043 TRICHLOROFUOROMETHANE
16	CI10 1,4-DIFLUOROBENZENE **IS#2**
17	C110 2-BUTANONE
18	C125 VINYL ACETATE
19	C120 CARBON TETRACHLORIDE
20	C130 BROMODICHLOROMETHANE
21	C140 1,2-DICHLOROPROPANE
22	C145 CIS-1,3-DICHLOROPROPENE
23	C150 TRICHLOROETHENE
24	C165 BENZENE
25	C155 DIBROMOCHLOROMETHANE
26	C160 1,1,2-TRICHLOROETHANE
27	C170 TRANS-1,3-DICHLOROPROPENE
28	C180 BROMOFORM
29	C115 1,1,1-TRICHLOROETHANE
30	CS05 TOLUENE-D8 **SS#2**
31	C120 CHLOROBENZENE-D5 **IS#3**
32	C205 4-METHYL-2-PENTANONE
33	C210 2-HEXANONE
34	C220 TETRACHLOROETHENE
35	C225 1,1,2,2-TETRACHLOROETHANE
36	C230 TOLUENE
37	C235 CHLOROBENZENE
38	C240 ETHYLBENZENE
39	C245 STYRENE
40	C250 ORTHO & PARA XYLENE
41	C251 META-XYLENE
42	CS10 BROMOFLUOROBENZENE **SS#3**

Ready 4D

100043

43394

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	128	254	10:35	1	1.000	A BB	381761	50.000 UG/L*	17.78
2	NOT FOUND								
3	NOT FOUND								
4	NOT FOUND								
5	NOT FOUND								
6	84	181	7:32	1	0.713	A BB	1005.	0.759 UG/L	0.34
7	NOT FOUND								
8	NOT FOUND								
9	NOT FOUND								
10	NOT FOUND								
11	NOT FOUND								
12	NOT FOUND								
13	NOT FOUND								
14	65	319	13:17	1	1.256	A BB	62189.	35.604 UG/L%	12.6671%
15	NOT FOUND								
16	114	498	20:45	16	1.000	A BB	164292.	50.000 UG/L*	17.78
17	NOT FOUND								
18	NOT FOUND								
19	NOT FOUND								
20	NOT FOUND								
21	NOT FOUND								
22	NOT FOUND								
23	NOT FOUND								
24	NOT FOUND								
25	NOT FOUND								
26	NOT FOUND								
27	NOT FOUND								
28	NOT FOUND								
29	NOT FOUND								
30	98	587	24:27	31	0.953	A BB	179639.	51.263 UG/L%	18.23
31	117	616	25:40	31	1.000	A BB	145738.	50.000 UG/L*	17.78
32	NOT FOUND								
33	NOT FOUND								
34	NOT FOUND								
35	NOT FOUND								
36	92	591	24:37	31	0.959	A BB	1063.	0.531 UG/L	0.19
37	NOT FOUND								
38	NOT FOUND								
39	NOT FOUND								
40	NOT FOUND								
41	NOT FOUND								
42	95	732	30:30	31	1.188	A BB	108505.	42.839 UG/L%	15.24

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
1	10:32	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
2	2:12		0.209						
3	3:22		0.320						
4	4:15		0.403						
5	5:15		0.498						
6	7:27	1.01	0.708	1.01	0.96	50.00	0.026	1.372	0.02
7	8:02		0.763						
8	9:02		0.858						
9	10:07		0.960						
10	11:25		1.083						
11	12:05		1.146						
12	12:40		1.202						

100044

43394

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
13	13:22		1. 269						
14	13:17	1. 00	1. 261	1. 00	35. 60	50. 00	1. 629	2. 288	0. 71
15	9:30		0. 901						
16	20:45	1. 00	1. 000	1. 00	50. 00	50. 00	1. 000	1. 000	1. 00
17	13:17		1. 261						
18	15:10		0. 731						
19	15:07		0. 729						
20	15:32		0. 749						
21	16:57		0. 817						
22	17:10		0. 827						
23	17:42		0. 853						
24	18:12		0. 878						
25	18:20		0. 884						
26	18:25		0. 888						
27	18:25		0. 888						
28	21:00		1. 012						
29	14:42		0. 709						
30	24:27	1. 00	0. 953	1. 00	51. 26	50. 00	1. 233	1. 202	1. 03
31	25:40	1. 00	1. 000	1. 00	50. 00	50. 00	1. 000	1. 000	1. 00
32	21:30		0. 838						
33	23:02		0. 898						
34	23:17		0. 907						
35	23:17		0. 907						
36	24:37	1. 00	0. 959	1. 00	0. 53	50. 00	0. 007	0. 686	0. 01
37	25:47		1. 005						
38	27:50		1. 084						
39	32:00		1. 247						
40	33:17		1. 297						
41	32:15		1. 256						
42	30:27	1. 00	1. 187	1. 00	42. 84	50. 00	0. 745	0. 869	0. 86

100045

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: VERSAR INC.

Contract: _____

5

Lab Code: VERSAR Case No.: R3-7

SAS No.: _____ SDG No.: 1

Matrix: (soil/water) SOIL

Lab Sample ID: 16429

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: U3413

Level: (low/med) LOW

Date Received: 04/19/90

% Moisture: not dec. 54

Date Analyzed: 04/27/90

Column: (pack/cap) PACK

Dilution Factor: 1.0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

Q

<u>74-87-3-----Chloromethane</u>	<u>22</u>	<u>U</u>
<u>74-83-9-----Bromomethane</u>	<u>22</u>	<u>U</u>
<u>75-01-4-----Vinyl chloride</u>	<u>22</u>	<u>U</u>
<u>75-00-3-----Chloroethane</u>	<u>22</u>	<u>U</u>
<u>75-09-2-----Methylene chloride</u>	<u>12</u>	
<u>67-64-1-----Acetone</u>	<u>32</u>	
<u>75-15-0-----Carbon disulfide</u>	<u>11</u>	<u>U</u>
<u>75-35-4-----1,1-Dichloroethene</u>	<u>11</u>	<u>U</u>
<u>75-34-3-----1,1-Dichloroethane</u>	<u>11</u>	<u>U</u>
<u>540-59-0-----1,2-Dichloroethene (total)</u>	<u>11</u>	<u>U</u>
<u>67-66-3-----Chloroform</u>	<u>11</u>	<u>U</u>
<u>107-06-2-----1,2-Dichloroethane</u>	<u>11</u>	<u>U</u>
<u>78-93-3-----2-Butanone</u>	<u>22</u>	<u>U</u>
<u>71-55-6-----1,1,1-Trichloroethane</u>	<u>11</u>	<u>U</u>
<u>56-23-5-----Carbon tetrachloride</u>	<u>11</u>	<u>U</u>
<u>108-05-4-----Vinyl acetate</u>	<u>22</u>	<u>U</u>
<u>75-27-4-----Bromodichloromethane</u>	<u>11</u>	<u>U</u>
<u>78-87-5-----1,2-Dichloropropane</u>	<u>11</u>	<u>U</u>
<u>10061-01-5-----cis-1,3-Dichloropropene</u>	<u>11</u>	<u>U</u>
<u>79-01-6-----Trichloroethene</u>	<u>11</u>	<u>U</u>
<u>124-48-1-----Dibromochloromethane</u>	<u>11</u>	<u>U</u>
<u>79-00-5-----1,1,2-Trichloroethane</u>	<u>11</u>	<u>U</u>
<u>71-43-2-----Benzene</u>	<u>11</u>	<u>U</u>
<u>10061-02-6-----Trans-1,3-dichloropropene</u>	<u>11</u>	<u>U</u>
<u>75-25-2-----Bromoform</u>	<u>11</u>	<u>U</u>
<u>108-10-1-----4-Methyl-2-pentanone</u>	<u>22</u>	<u>U</u>
<u>591-78-6-----2-Hexanone</u>	<u>22</u>	<u>U</u>
<u>127-18-4-----Tetrachloroethene</u>	<u>11</u>	<u>U</u>
<u>79-34-5-----1,1,2,2-Tetrachloroethane</u>	<u>11</u>	<u>U</u>
<u>108-88-3-----Toluene</u>	<u>11</u>	<u>U</u>
<u>108-90-7-----Chlorobenzene</u>	<u>11</u>	<u>U</u>
<u>100-41-4-----Ethylbenzene</u>	<u>11</u>	<u>U</u>
<u>100-42-5-----Styrene</u>	<u>11</u>	<u>U</u>
<u>1330-20-7-----Total xylenes</u>	<u>11</u>	<u>U</u>

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: VERSAR INC.

Contract: _____

5

Lab Code: VERSAR Case No.: R3-7

SAS No.: _____ SDG No.: 1 _____

Matrix: (soil/water) SOIL

Lab Sample ID: 16429 _____

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: U3413 _____

Level: (low/med) LOW

Date Received: 04/19/90

% Moisture: not dec. 54

Date Analyzed: 04/27/90

Column (pack/cap) PACK

Dilution Factor: 1.0

Number TICs found: 0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

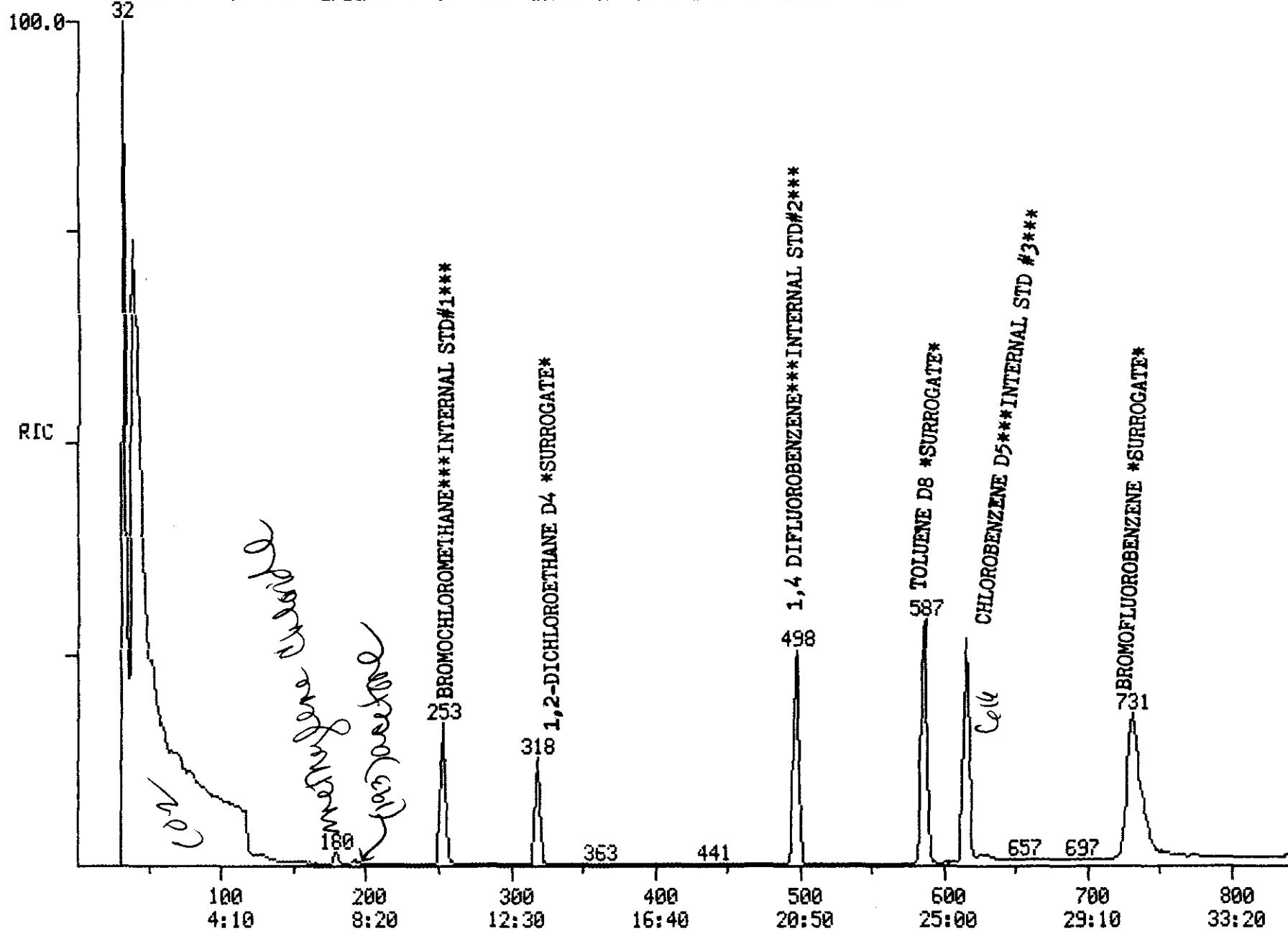
RIC
04/27/90 10:22:00

SAMPLE: CLP, VERSCDM, 2536, 85, L, S, 16429, V, .420, 1.0, B2, 5ML,,

COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

RANGE: G 1, 840 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: U3413 #1 SCANS 1 TO 840
CALI: U3413 #2 55/Surrogate



Quantitation Report File: U3413

Data: U3413.TI

04/27/90 10:22:00

Sample: CLP, VERSCDM, 2536, 85, L, S, 16429, V, , 420.1.0, B2, 5ML

Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

Formula: -

Instrument: U

Weight: 0.011

Submitted by: VERSAR

Analyst: MAT

Acct. No.: 420.1.0

5g/5ml 1/1

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)

Resp. fac. from Library Entry

54.5 %

No	Name
1	C101 BROMOCHLOROMETHANE **IS#1**
2	C010 CHLOROMETHANE
3	C015 BROMOMETHANE
4	C020 VINYL CHLORIDE
5	C025 CHLOROETHANE
6	C030 METHYLENE CHLORIDE
7	C035 ACETONE
8	C040 CARBON DISULFIDE
9	C045 1,1-DICHLOROETHENE
10	C050 1,1-DICHLOROETHANE
11	C053 1,2-DICHLOROETHENE (TOTAL)
12	C060 CHLOROFORM
13	C065 1,2-DICHLOROETHANE
14	CS15 1,2-DICHLOROETHANE-D4 **SS#1**
15	C043 TRICHLOROFUOROMETHANE
16	CI10 1,4-DIFLUOROBENZENE **IS#2**
17	C110 2-BUTANONE
18	C125 VINYL ACETATE
19	C120 CARBON TETRACHLORIDE
20	C130 BROMODICHLOROMETHANE
21	C140 1,2-DICHLOROPROPANE
22	C145 CIS-1,3-DICHLOROPROPENE
23	C150 TRICHLOROETHENE
24	C165 BENZENE
25	C155 DIBROMOCHLOROMETHANE
26	C160 1,1,2-TRICHLOROETHANE
27	C170 TRANS-1,3-DICHLOROPROPENE
28	C180 BROMOFORM
29	C115 1,1,1-TRICHLOROETHANE
30	CS05 TOLUENE-D8 **SS#2**
31	CI20 CHLOROBENZENE-D5 **IS#3**
32	C205 4-METHYL-2-PENTANONE
33	C210 2-HEXANONE
34	C220 TETRACHLOROETHENE
35	C225 1,1,2,2-TETRACHLOROETHANE
36	C230 TOLUENE
37	C235 CHLOROBENZENE
38	C240 ETHYLBENZENE
39	C245 STYRENE
40	C250 ORTHO & PARA XYLENE
41	C251 META-XYLENE
42	CS10 BROMOFLUOROBENZENE **SS#3**

INT STD & SURRG. COMPLIANT

Ready to

10049

U3413

100050

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	128	253	10:32	1	1.000	A BB	58714. ✓	50.000 UG/L*	15.89
2	NOT FOUND								
3	NOT FOUND								
4	NOT FOUND								
5	NOT FOUND								
6	84	180	7:30	1	0.711	A BB	8066.	5.415 UG/L	1.72
7	43	193	8:02	1	0.763	A BV	8540.	14.712 UG/L	4.67
8	NOT FOUND								
9	NOT FOUND								
10	NOT FOUND								
11	NOT FOUND								
12	NOT FOUND								
13	NOT FOUND								
14	65	318	13:15	1	1.257	A BB	98321.	41.690 UG/L%	13.25 83%
15	NOT FOUND								
16	114	498	20:45	16	1.000	A BB	241997. ✓	50.000 UG/L*	15.89
17	72	318	13:15	1	1.257	A BB	179.	0.772 UG/L	0.25
18	NOT FOUND								
19	NOT FOUND								
20	NOT FOUND								
21	NOT FOUND								
22	NOT FOUND								
23	NOT FOUND								
24	78	437	18:12	16	0.878	A BB	733.	0.168 UG/L	0.05
25	NOT FOUND								
26	NOT FOUND								
27	NOT FOUND								
28	NOT FOUND								
29	NOT FOUND								
30	98	586	24:25	31	0.951	A BB	249604.	54.071 UG/L%	17.18 108%
31	117	616	25:40	31	1.000	A BB	200802. ✓	50.000 UG/L*	15.89
32	43	516	21:30	31	0.838	A BB	767.	0.264 UG/L	0.08
33	NOT FOUND								
34	NOT FOUND								
35	NOT FOUND								
36	92	591	24:37	31	0.959	A BB	2495.	0.974 UG/L	0.31
37	NOT FOUND								
38	NOT FOUND								
39	NOT FOUND								
40	NOT FOUND								
41	106	774	32:15	31	1.256	A BB	683.	0.337 UG/L	0.11
42	95	731	30:27	31	1.187	A BB	140128.	46.324 UG/L%	14.72 93%

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
1	10:30	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
2	2:10		0.206						
3	3:20		0.317						
4	4:12		0.401						
5	5:15		0.500						
6	7:25	1.01	0.706	1.01	5.42	50.00	0.137	1.268	0.11
7	8:00	1.01	0.762	1.00	14.71	50.00	0.145	0.494	0.29
8	9:00		0.857						
9	10:05		0.960						
10	11:22		1.083						
11	12:02		1.147						
12	12:37		1.202						

100050

43413

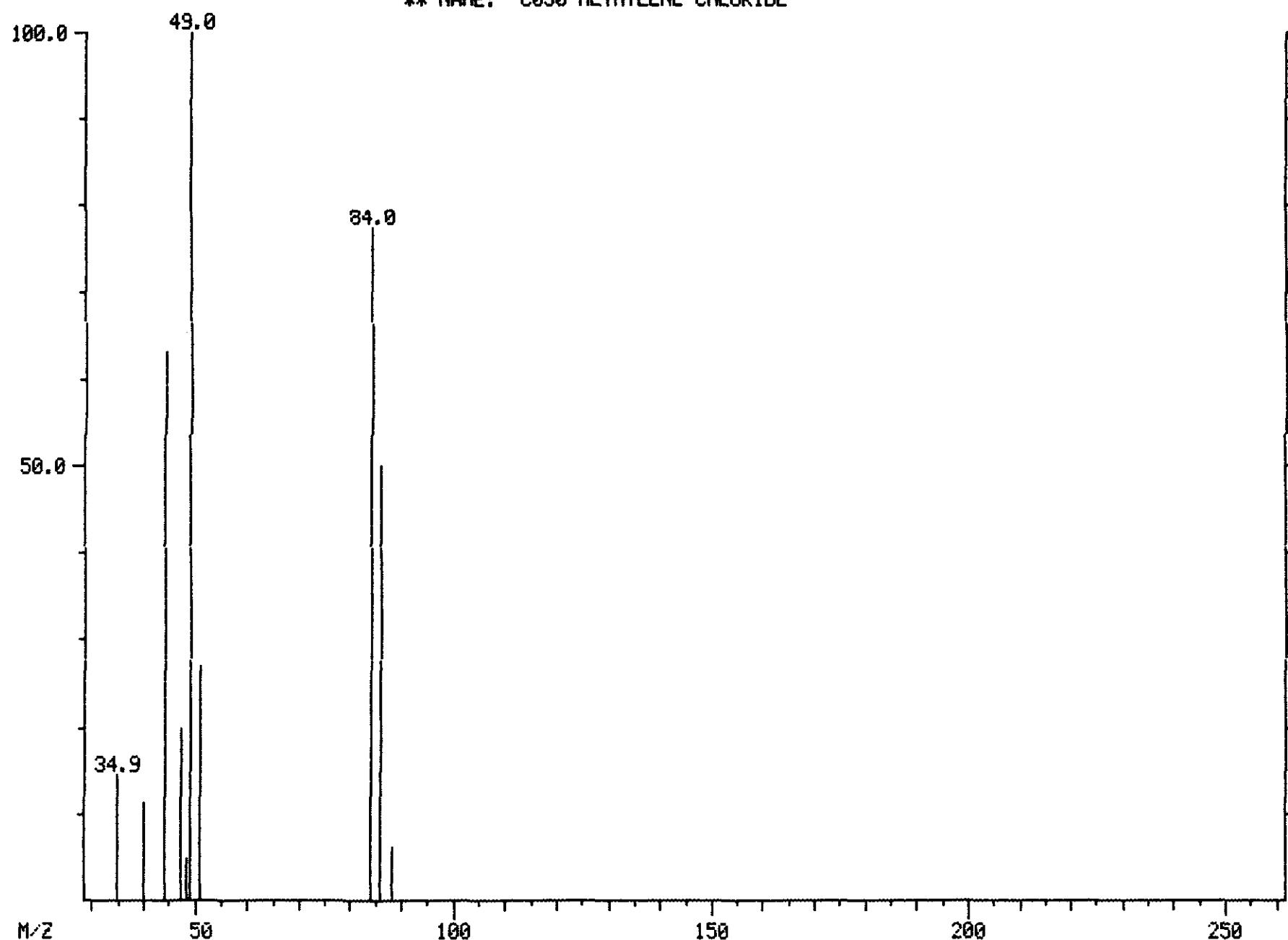
No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R.Fac	R.Fac(L)	Ratio
13	13: 20		1. 270						
14	13: 15	1. 00	1. 262	1. 00	41. 69	50. 00	1. 675	2. 008	0. 83
15	9: 27		0. 901						
16	20: 42	1. 00	1. 000	1. 00	50. 00	50. 00	1. 000	1. 000	1. 00
17	13: 15	1. 00	1. 262	1. 00	0. 77	50. 00	0. 003	0. 197	0. 02
18	15: 07		0. 730						
19	15: 05		0. 728						
20	15: 30		0. 748						
21	16: 55		0. 817						
22	17: 07		0. 827						
23	17: 40		0. 853						
24	18: 12	1. 00	0. 879	1. 00	0. 17	50. 00	0. 003	0. 901	0. 00
25	18: 17		0. 883						
26	18: 25		0. 889						
27	18: 22		0. 887						
28	20: 57		1. 012						
29	14: 42		0. 710						
30	24: 27	1. 00	0. 953	1. 00	54. 07	50. 00	1. 243	1. 149	1. 08
31	25: 40	1. 00	1. 000	1. 00	50. 00	50. 00	1. 000	1. 000	1. 00
32	21: 30	1. 00	0. 838	1. 00	0. 26	50. 00	0. 004	0. 725	0. 01
33	23: 02		0. 898						
34	23: 17		0. 907						
35	23: 17		0. 907						
36	24: 37	1. 00	0. 959	1. 00	0. 97	50. 00	0. 012	0. 638	0. 02
37	25: 47		1. 005						
38	27: 57		1. 089						
39	32: 25		1. 263						
40	33: 47		1. 317						
41	32: 42	0. 99	1. 274	0. 99	0. 34	50. 00	0. 003	0. 505	0. 01
42	30: 45	0. 99	1. 198	0. 99	46. 32	50. 00	0. 698	0. 753	0. 93

100051

MASS SPECTRUM
04/27/90 10:22:00 + 7:30
SAMPLE: CLP,VERSCDM,2536,B5,L,S,16429,U,,420.1.0,B2,5ML,,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
** NAME: C030 METHYLENE CHLORIDE

DATA: U3413 #180
CALI: U3413 #2
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
** NAME: C030 METHYLENE CHLORIDE

BASE M/Z: 49
RIC: 8768.



2344.1 00052

MASS SPECTRUM

04/27/90 10:22:00 + 7:30

SAMPLE: CLP,VERSCDM,2536,B5,L,S,16429,V,,420.1.0,B2,5ML,,

COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

** NAME: C030 METHYLENE CHLORIDE

ENHANCED (S 15B 2N 0T)

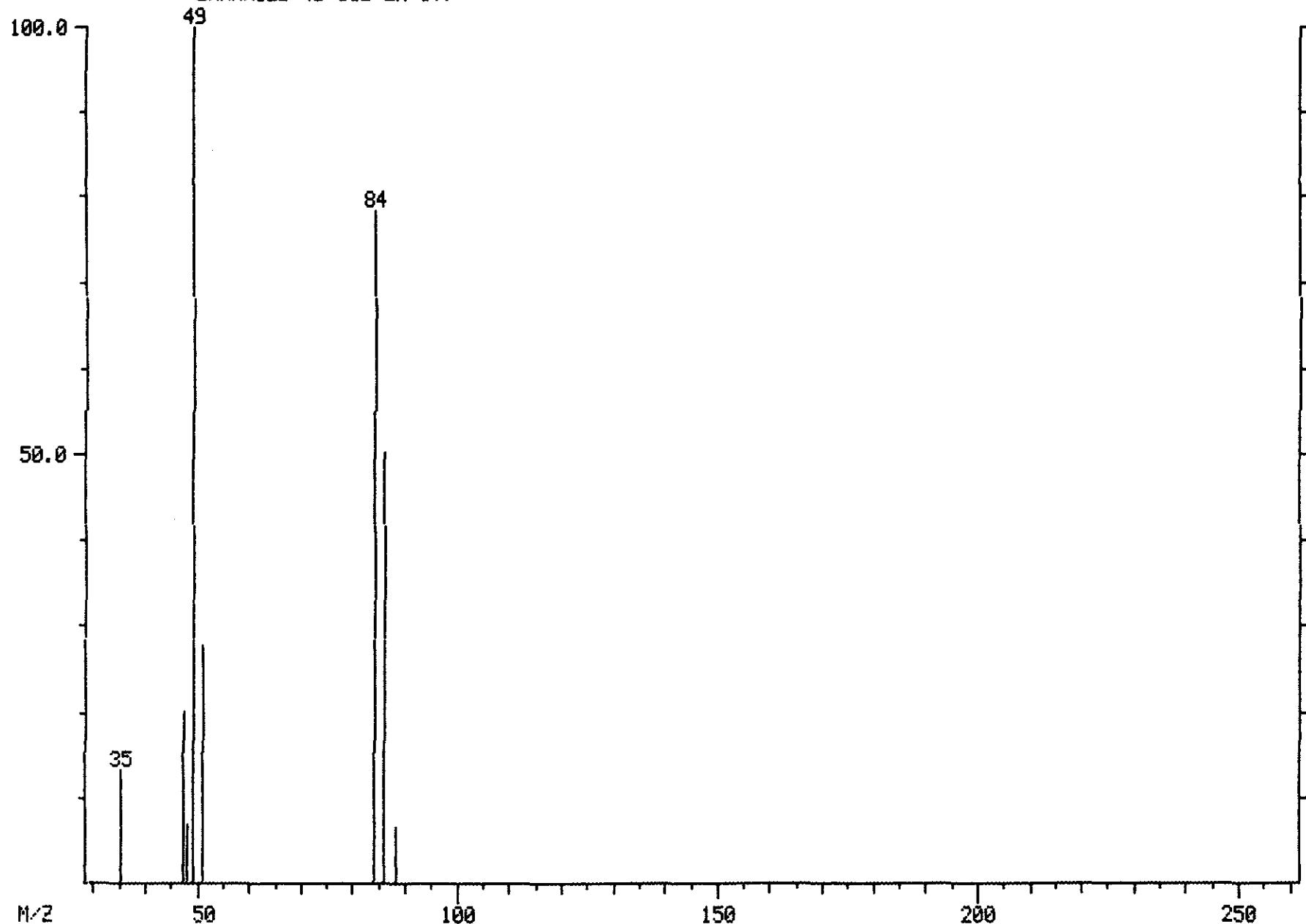
DATA: U3413 #180

CALI: U3413 #2

BASE M/Z: 49

RIC: 6896.

1100053
2268.



MASS SPECTRUM

04/27/90 8:21:00 + 7:25

SAMPLE: CLP,,,USTD 50,L,S,22554,U,CC-050,,5ML,,

COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

** NAME: C030 METHYLENE CHLORIDE

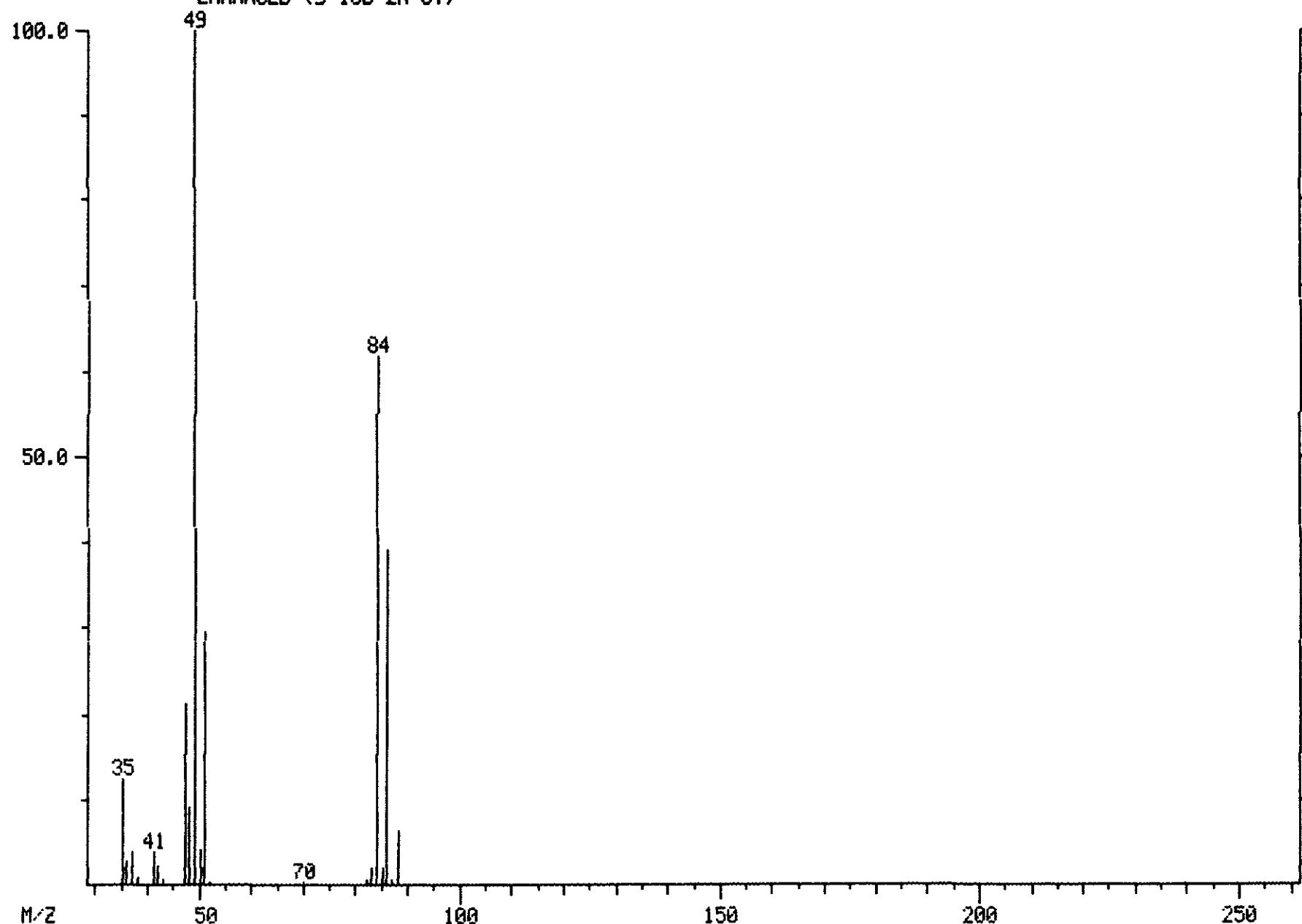
ENHANCED (S 15B 2N 0T)

DATA: U3411 #178

CALI: U3411 #2

BASE M/Z: 49

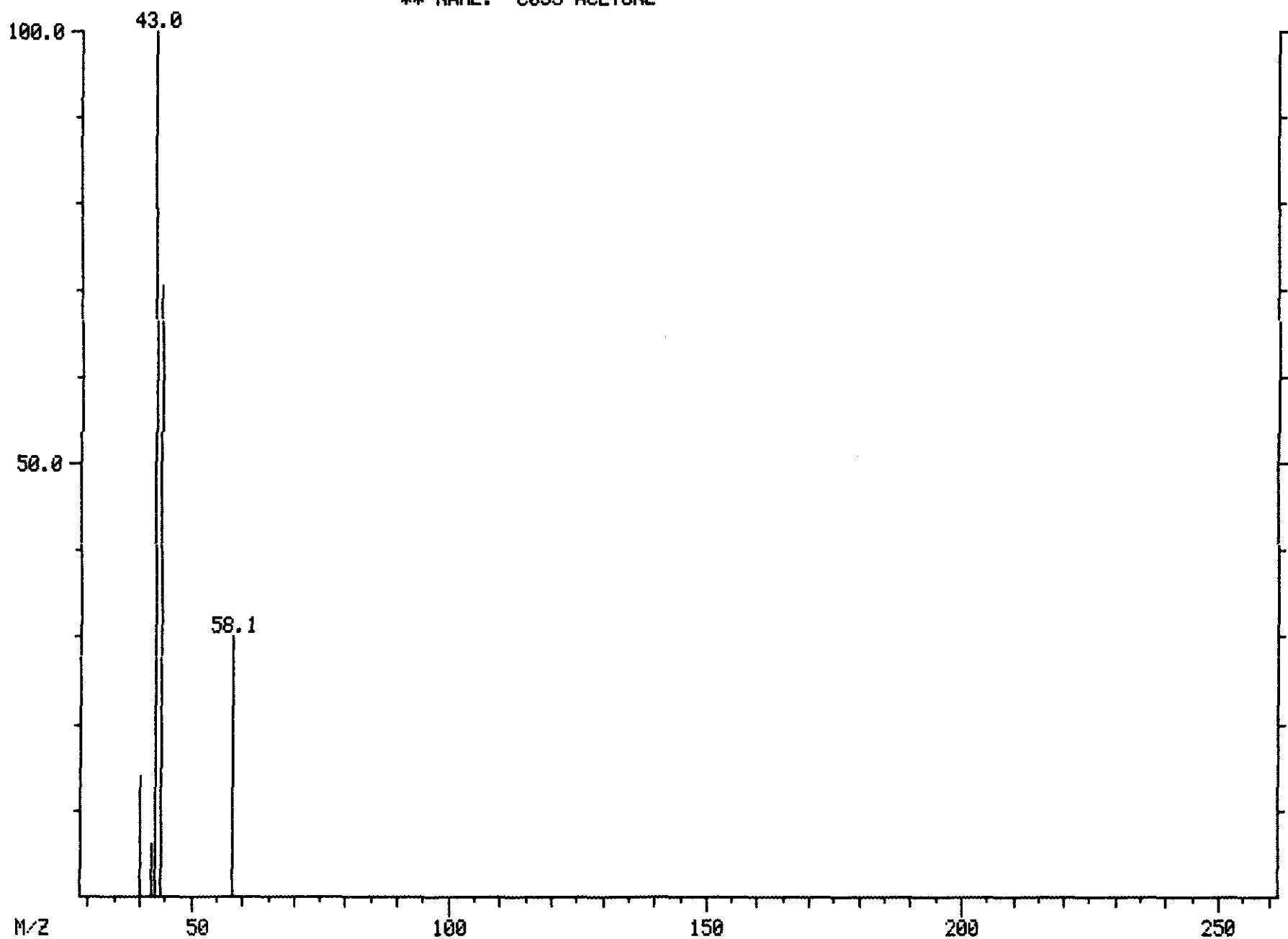
RIC: 103936.



1000054

MASS SPECTRUM
04/27/90 10:22:00 + 8:02
SAMPLE: CLP,VERSCDM,2536,B5,L,S,16429,U,,420.1.0,B2,5ML,,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
** NAME: C035 ACETONE

DATA: U3413 #193
CALI: U3413 #2
BASE M/Z: 43
RIC: 4184.

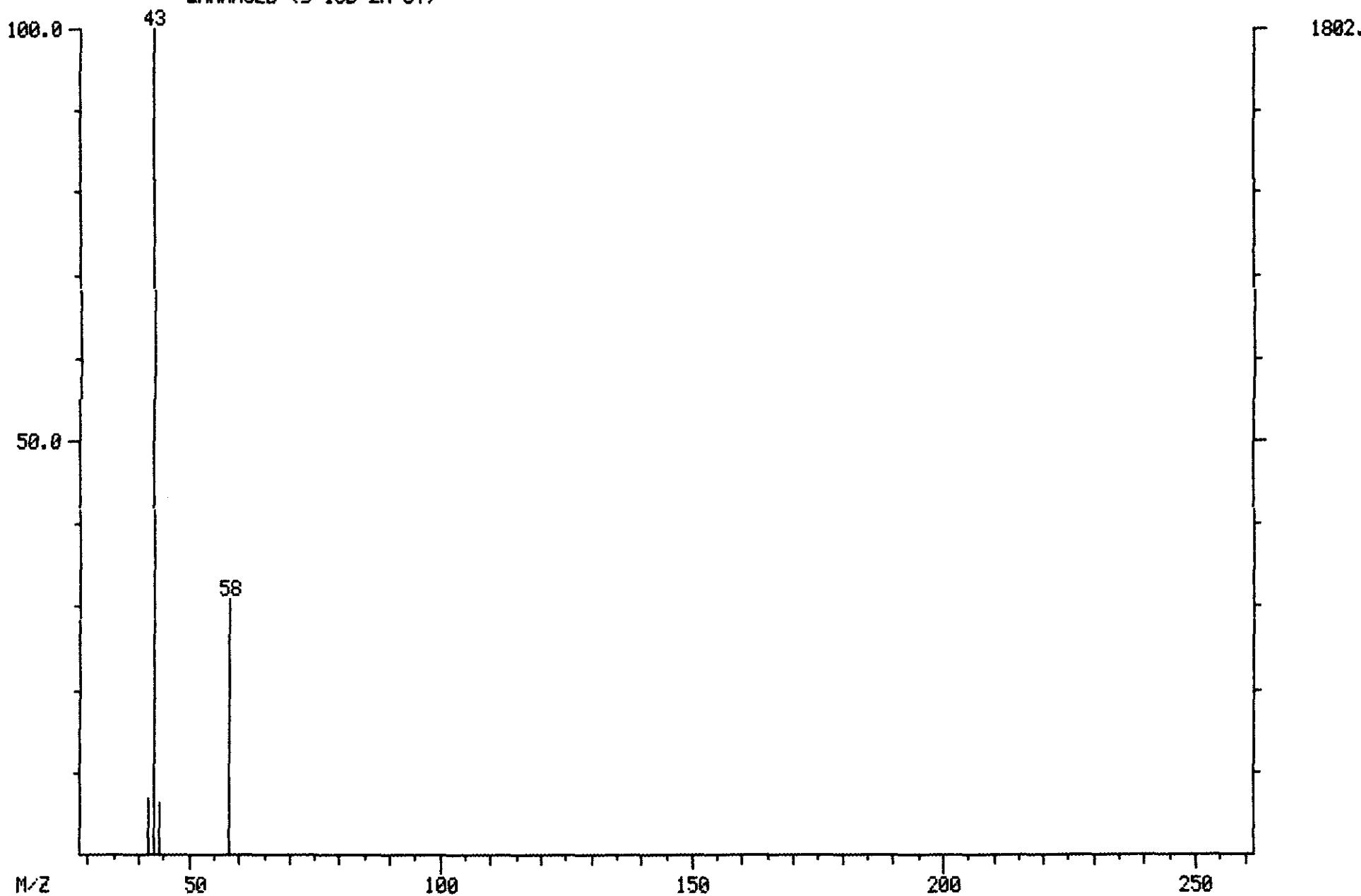


00055
1898.1

MASS SPECTRUM
04/27/90 10:22:00 + 8:02
SAMPLE: CLP,VER5CDM,2536,B5,L,S,16429,U,,420.1.0,B2,5ML,,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
** NAME: C035 ACETONE
ENHANCED (S 15B 2N 0T)

DATA: U3413 #193
CALI: U3413 #2

BASE M/Z: 43
RIC: 2604.



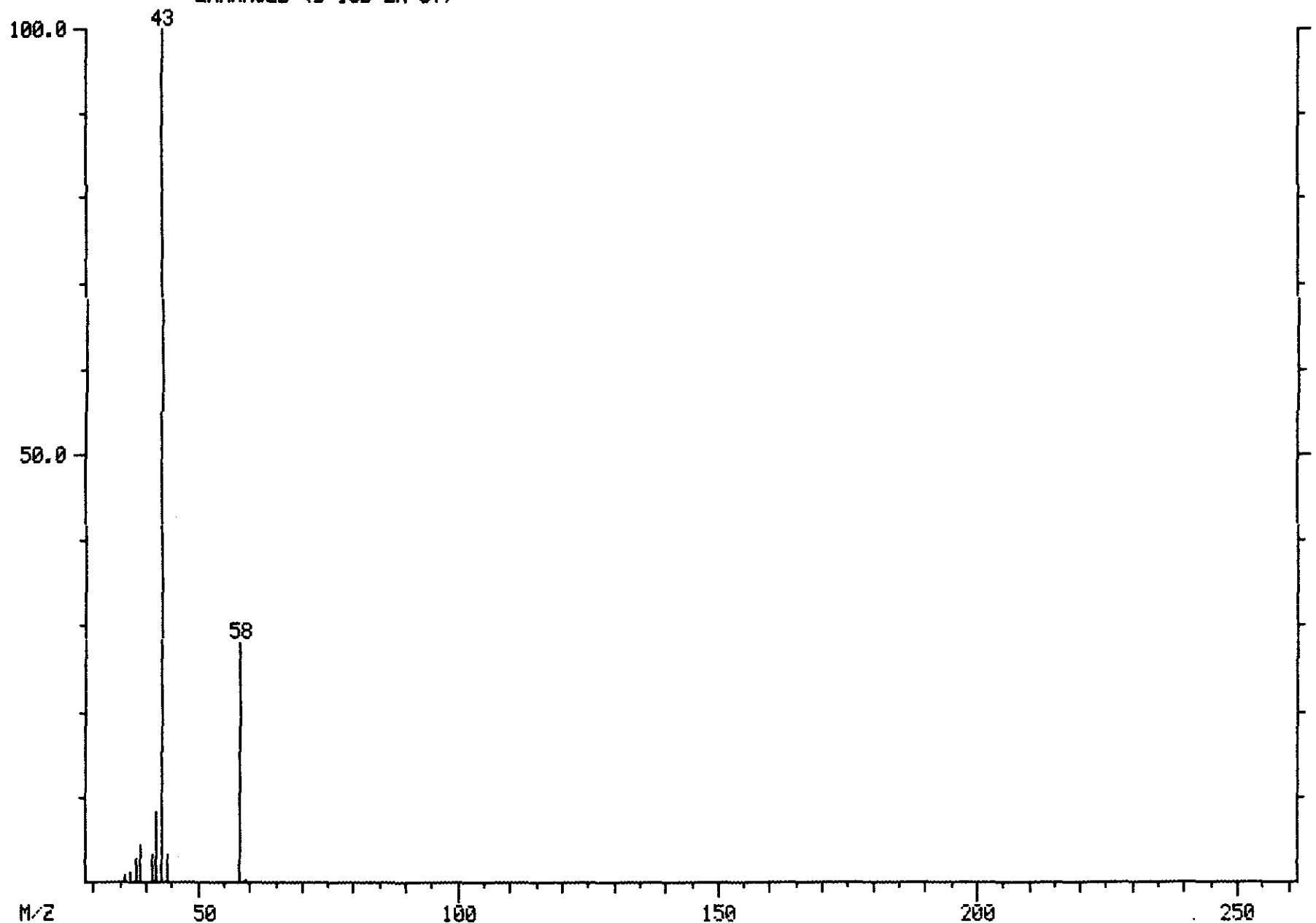
100056

MASS SPECTRUM
04/27/90 8:21:00 + 8:00
SAMPLE: CLP,,,VSTD 50,L,S,22554,U,CC-050,,5ML,,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
** NAME: C835 ACETONE
ENHANCED (S 15B 2N 0T)

DATA: U3411 #192
CALI: U3411 #2
CONDS.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
** NAME: C835 ACETONE
ENHANCED (S 15B 2N 0T)

BASE M/Z: 43
RIC: 12368.

1000057



1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: VERSAR INC.

Contract: _____

6

Lab Code: VERSAR Case No.: R3-7

SAS No.: _____ SDG No.: 1

Matrix: (soil/water) SOIL

Lab Sample ID: 16430

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: U3414

Level: (low/med) LOW

Date Received: 04/19/90

% Moisture: not dec. 40

Date Analyzed: 04/27/90

Column: (pack/cap) PACK

Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3-----	Chloromethane	17	U
74-83-9-----	Bromomethane	17	U
75-01-4-----	Vinyl chloride	17	U
75-00-3-----	Chloroethane	17	U
75-09-2-----	Methylene chloride	8	U
67-64-1-----	Acetone	31	
75-15-0-----	Carbon disulfide	8	U
75-35-4-----	1,1-Dichloroethene	8	U
75-34-3-----	1,1-Dichloroethane	8	U
540-59-0-----	1,2-Dichloroethene (total)	8	U
67-66-3-----	Chloroform	8	U
107-06-2-----	1,2-Dichloroethane	8	U
78-93-3-----	2-Butanone	17	U
71-55-6-----	1,1,1-Trichloroethane	8	U
56-23-5-----	Carbon tetrachloride	8	U
108-05-4-----	Vinyl acetate	17	U
75-27-4-----	Bromodichloromethane	8	U
78-87-5-----	1,2-Dichloropropane	8	U
10061-01-5-----	cis-1,3-Dichloropropene	8	U
79-01-6-----	Trichloroethene	8	U
124-48-1-----	Dibromochloromethane	8	U
79-00-5-----	1,1,2-Trichloroethane	8	U
71-43-2-----	Benzene	8	U
10061-02-6-----	Trans-1,3-dichloropropene	8	U
75-25-2-----	Bromoform	8	U
108-10-1-----	4-Methyl-2-pentanone	17	U
591-78-6-----	2-Hexanone	17	U
127-18-4-----	Tetrachloroethene	8	U
79-34-5-----	1,1,2,2-Tetrachloroethane	8	U
108-88-3-----	Toluene	8	U
108-90-7-----	Chlorobenzene	8	U
100-41-4-----	Ethylbenzene	8	U
100-42-5-----	Styrene	8	U
1330-20-7-----	Total xylenes	8	U

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: VERSAR INC.

Contract: _____

6

Lab Code: VERSAR Case No.: R3-7 SAS No.: _____ SDG No.: 1 _____

Matrix: (soil/water) SOIL

Lab Sample ID: 16430 _____

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: U3414 _____

Level: (low/med) LOW

Date Received: 04/19/90

* Moisture: not dec. 40

Date Analyzed: 04/27/90

Column (pack/cap) PACK

Dilution Factor: 1.0

Number TICs found: 0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

RIC
04/27/90 11:05:00

DATA: U3414 #1
CALI: U3414 #2

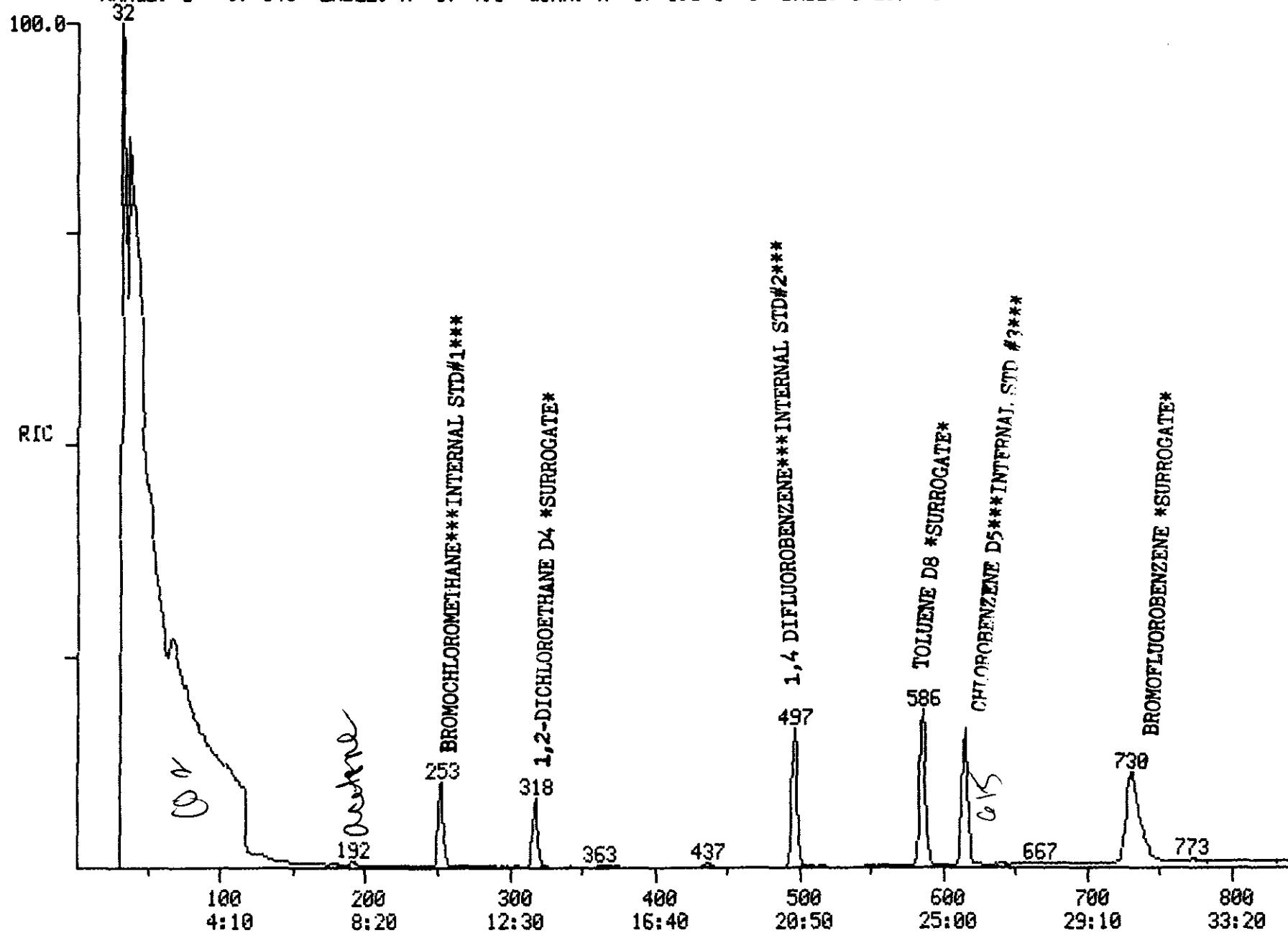
SCANS 1 TO 840

SAMPLE: CLP, VERSCOM, 2536, B6, L, S, 16430, U, , 420.1.0, B2, 5G/5ML,

COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

RANGE: G 1, 840 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

777216.
77100060



Quantitation Report File: U3414

Data: U3414.TI

04/27/90 11:05:00

Sample: CLP, VERSCDM, 2536, B6, L, S, 16430, V, , 420.1.0, B2, 5G/5ML, ,

Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

Formula: -

Instrument: U

Weight: 0.011

Submitted by: VERSAR

Analyst: MAT

Acct. No.: 420.1.0

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)

Resp. fac. from Library Entry

40.9 %

No	Name
1	C101 BROMOCHLOROMETHANE **IS#1**
2	C010 CHLOROMETHANE
3	C015 BROMOMETHANE
4	C020 VINYL CHLORIDE
5	C025 CHLOROETHANE
6	C030 METHYLENE CHLORIDE
7	C035 ACETONE
8	C040 CARBON DISULFIDE
9	C045 1,1-DICHLOROETHENE
10	C050 1,1-DICHLOROETHANE
11	C053 1,2-DICHLOROETHENE (TOTAL)
12	C060 CHLOROFORM
13	C065 1,2-DICHLOROETHANE
14	CS15 1,2-DICHLOROETHANE-D4 **SS#1**
15	C043 TRICHLOROFUOROMETHANE
16	CI10 1,4-DIFLUOROBENZENE **IS#2**
17	C110 2-BUTANONE
18	C125 VINYL ACETATE
19	C120 CARBON TETRACHLORIDE
20	C130 BROMODICHLOROMETHANE
21	C140 1,2-DICLOROPROPANE
22	C145 CIS-1,3-DICHLOROPROPENE
23	C150 TRICHLOROETHENE
24	C165 BENZENE
25	C155 DIBROMOCHLOROMETHANE
26	C160 1,1,2-TRICHLOROETHANE
27	C170 TRANS-1,3-DICHLOROPROPENE
28	C180 BROMOFORM
29	C115 1,1,1-TRICHLOROETHANE
30	CS05 TOLUENE-D8 **SS#2**
31	CI20 CHLOROBENZENE-D5 **IS#3**
32	C205 4-METHYL-2-PENTANONE
33	C210 2-HEXANONE
34	C220 TETRACHLOROETHENE
35	C225 1,1,2,2-TETRACHLOROETHANE
36	C230 TOLUENE
37	C235 CHLOROBENZENE
38	C240 ETHYLBENZENE
39	C245 STYRENE
40	C250 ORTHO & PARA XYLENE
41	C251 META-XYLENE
42	CS10 BROMOFLUOROBENZENE **SS#3**

Realty

100061

W3414

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	128	252	10:30	1	1.000	A BB	53154. ✓	50.000 UG/L*	15.39
2	NOT FOUND								
3	NOT FOUND								
4	NOT FOUND								
5	NOT FOUND								
6	84	179	7:27	1	0.710	A BB	2727.	12.022 UG/L	0.62
7	43	192	8:00	1	0.762	A BB	9800.	<u>18.649 UG/L</u>	5.74
8	NOT FOUND								
9	NOT FOUND								
10	63	253	10:32	1	1.004	A BB	176.	0.068 UG/L	0.02
11	NOT FOUND								
12	NOT FOUND								
13	NOT FOUND								
14	65	318	13:15	1	1.262	A BB	89424.	41.884 UG/L%	12.89 84%
15	NOT FOUND								
16	114	497	20:42	16	1.000	A BB	226023. ✓	50.000 UG/L*	15.39
17	72	319	13:17	1	1.266	A BB	963.	4.593 UG/L	1.41
18	NOT FOUND								
19	NOT FOUND								
20	NOT FOUND								
21	NOT FOUND								
22	NOT FOUND								
23	NOT FOUND								
24	78	437	18:12	16	0.879	A BB	7700.	1.890 UG/L	0.58
25	NOT FOUND								
26	NOT FOUND								
27	NOT FOUND								
28	NOT FOUND								
29	NOT FOUND								
30	98	586	24:25	31	0.953	A BB	230466.	55.302 UG/L%	17.03 11%
31	117	615	25:37	31	1.000	A BB	181280.✓	50.000 UG/L*	15.39
32	NOT FOUND								
33	NOT FOUND								
34	NOT FOUND								
35	NOT FOUND								
36	NOT FOUND								
37	NOT FOUND								
38	NOT FOUND								
39	NOT FOUND								
40	106	774	32:15	31	1.259	A BB	3078.	1.940 UG/L	0.60
41	106	774	32:15	31	1.259	A BB	3078.	1.681 UG/L	0.52
42	95	731	30:27	31	1.189	A BB	127832.	46.810 UG/L%	14.41 94%

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
1	10:30	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
2	2:10		0.206						
3	3:20		0.317						
4	4:12		0.401						
5	5:15		0.500						
6	7:25	1.01	0.706	1.01	2.02	50.00	0.051	1.268	0.04
7	8:00	1.00	0.762	1.00	18.65	50.00	0.184	0.494	0.37
8	9:00		0.857						
9	10:05		0.960						
10	11:22	0.93	1.083	0.93	0.07	50.00	0.003	2.426	0.00
11	12:02		1.147						
12	12:37		1.202						

100062

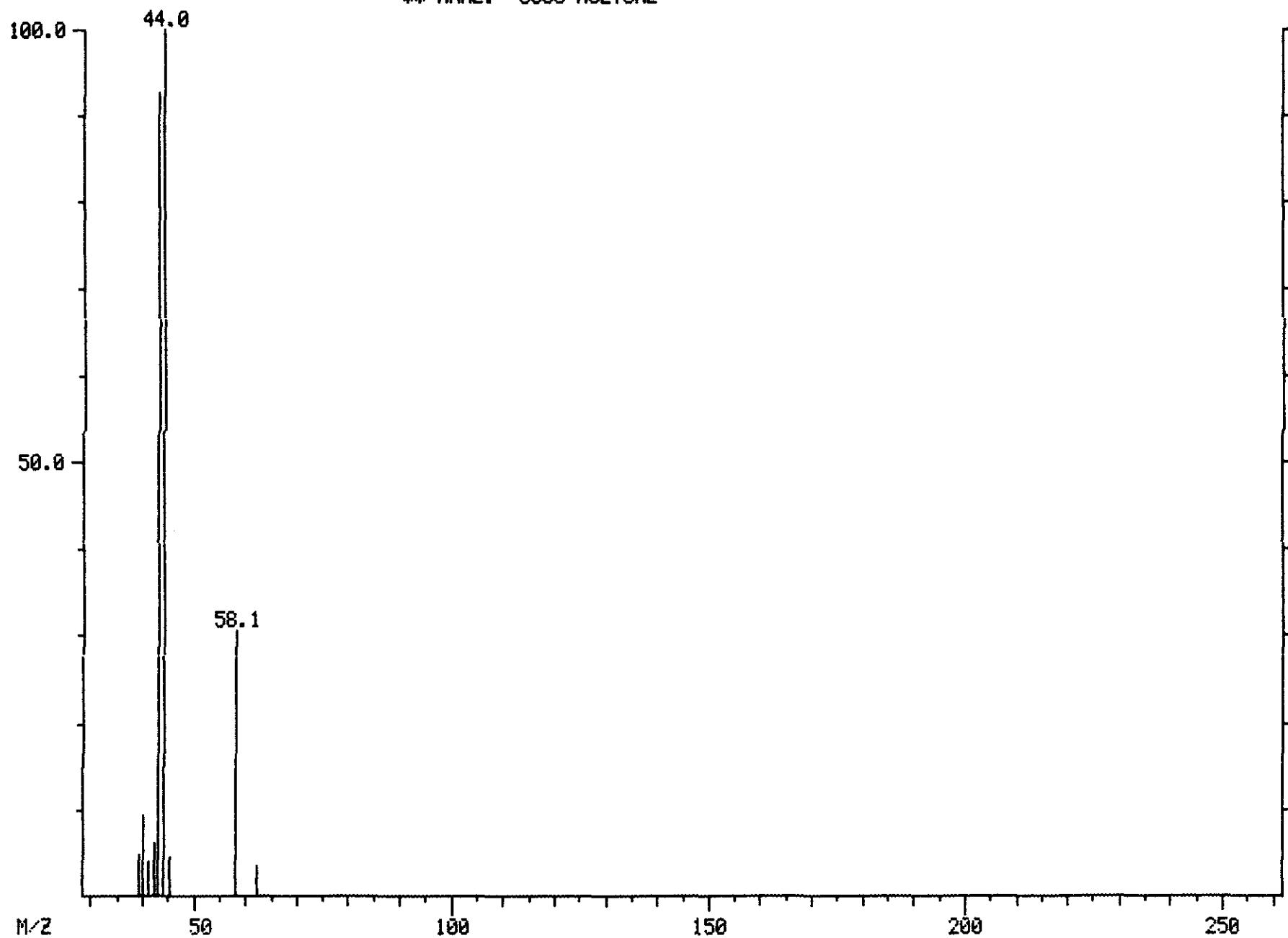
u3414

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
13	13:20		1. 270						
14	13:15	1. 00	1. 262	1. 00	41. 88	50. 00	1. 682	2. 008	0. 84
15	9:27		0. 901						
16	20:42	1. 00	1. 000	1. 00	50. 00	50. 00	1. 000	1. 000	1. 00
17	13:15	1. 00	1. 262	1. 00	4. 59	50. 00	0. 018	0. 197	0. 09
18	15:07		0. 730						
19	15:05		0. 728						
20	15:30		0. 748						
21	16:55		0. 817						
22	17:07		0. 827						
23	17:40		0. 853						
24	18:12	1. 00	0. 879	1. 00	1. 89	50. 00	0. 034	0. 901	0. 04
25	18:17		0. 883						
26	18:25		0. 889						
27	18:22		0. 887						
28	20:57		1. 012						
29	14:42		0. 710						
30	24:27	1. 00	0. 953	1. 00	55. 30	50. 00	1. 271	1. 149	1. 11
31	25:40	1. 00	1. 000	1. 00	50. 00	50. 00	1. 000	1. 000	1. 00
32	21:30		0. 838						
33	23:02		0. 898						
34	23:17		0. 907						
35	23:17		0. 907						
36	24:37		0. 959						
37	25:47		1. 005						
38	27:57		1. 089						
39	32:25		1. 263						
40	33:47	0. 95	1. 317	0. 96	1. 94	50. 00	0. 017	0. 438	0. 04
41	32:42	0. 99	1. 274	0. 99	1. 68	50. 00	0. 017	0. 505	0. 03
42	30:45	0. 99	1. 198	0. 99	46. 81	50. 00	0. 705	0. 753	0. 94

MASS SPECTRUM
04/27/90 11:05:00 + 8:00
SAMPLE: CLP,VERSCDM,2536,B6,L,S,16430,U,,420.1.0,B2,5G/5ML,,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
** NAME: C035 ACETONE

DATA: U3414 #192
CALI: U3414 #2
BASE M/Z: 44
RIC: 6080.

1000064
2372.1



MASS SPECTRUM

04/27/90 11:05:00 + 8:00

SAMPLE: CLP,VERSCDM,2536,B6,L,S,16430,U,,420.1.0,B2,5G/5ML,,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
** NAME: C035 ACETONE

DATA: U3414 #192

CALI: U3414 #2

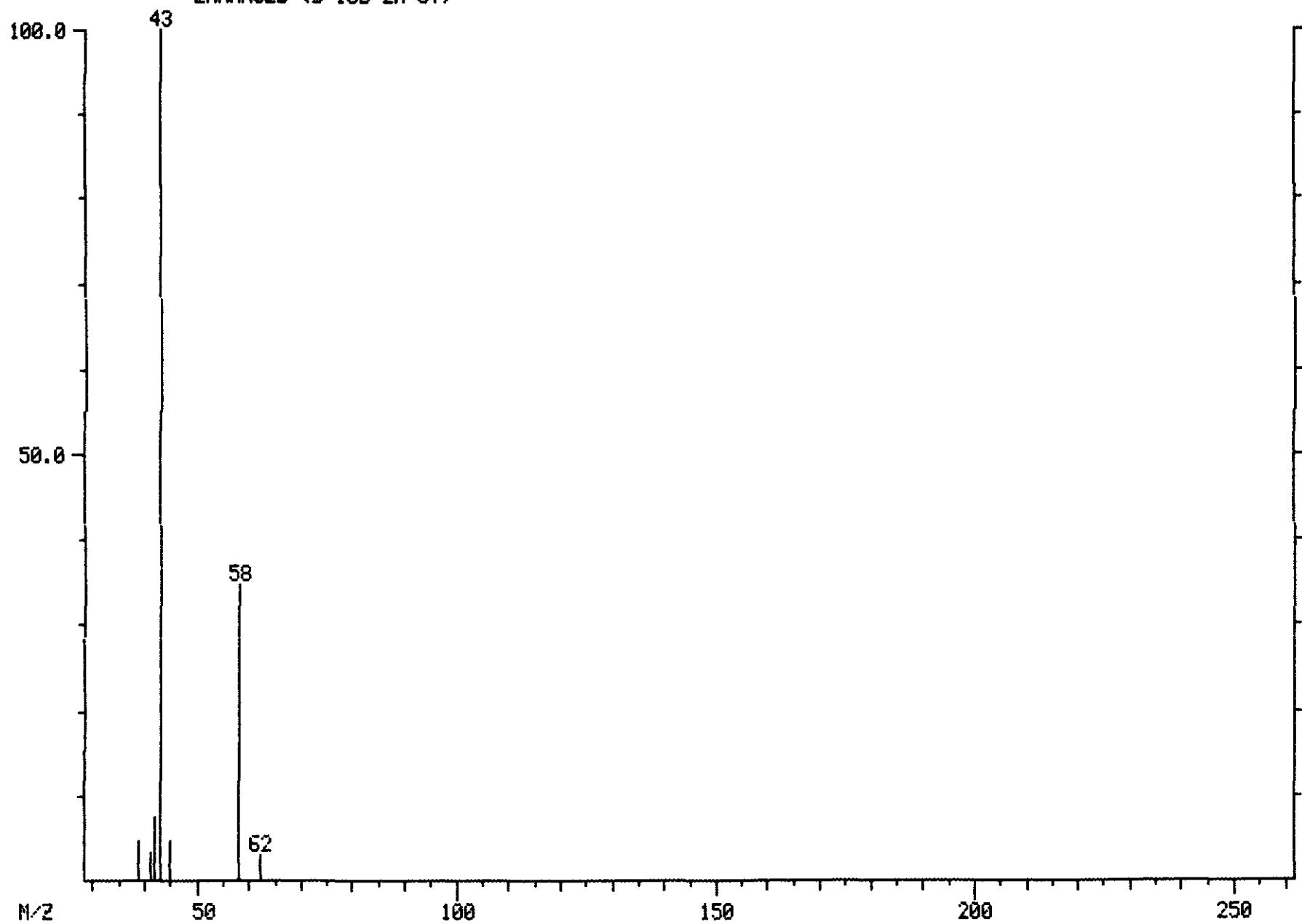
BASE M/Z: 43

RIC: 3352.

1000065

2120.

ENHANCED (S 15B 2N 0T)



MASS SPECTRUM

04/27/90 8:21:00 + 8:00

SAMPLE: CLP,,,VSTD 50,L,S,22554,U,CC-050,,5ML,,

COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

** NAME: C035 ACETONE

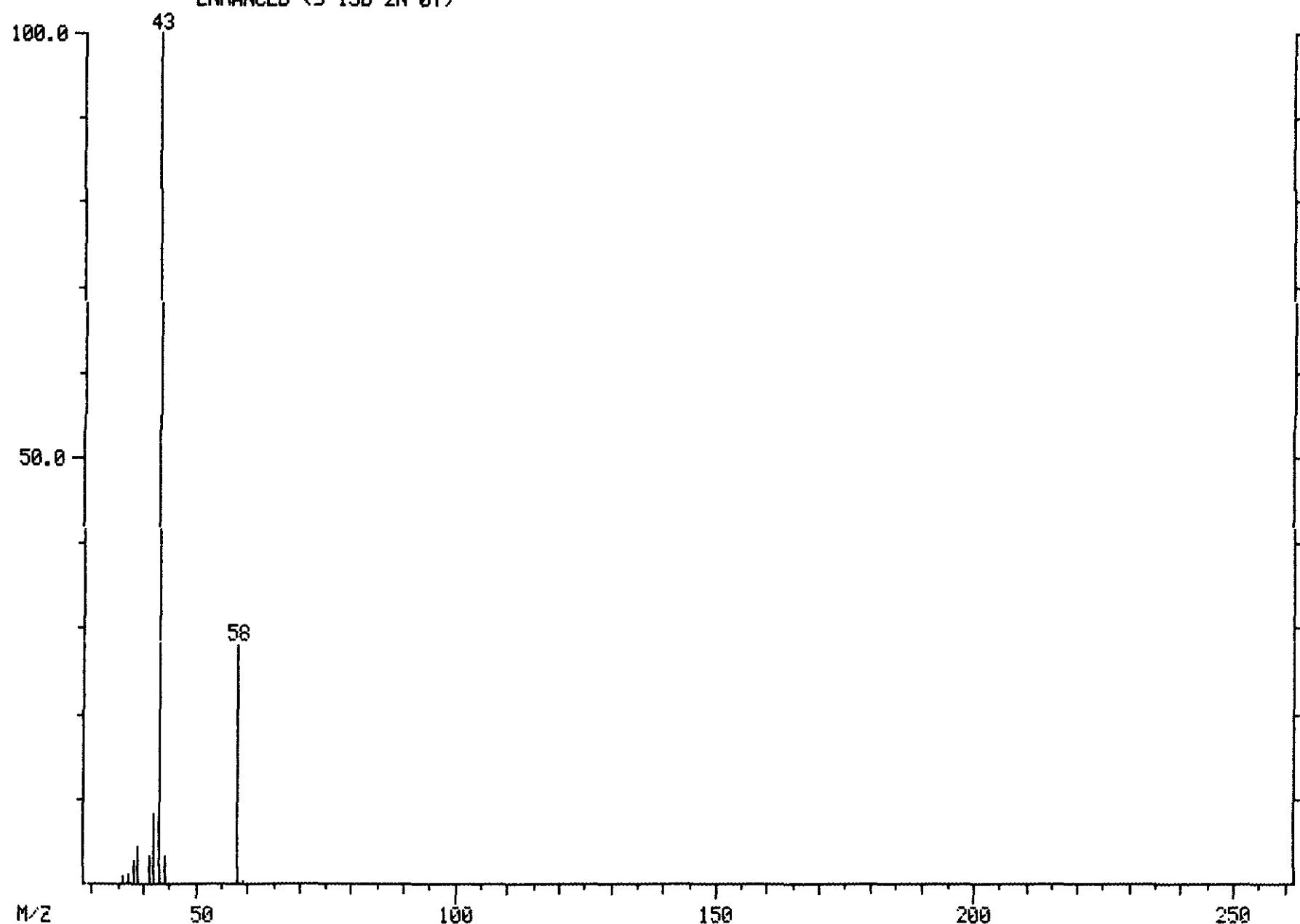
ENHANCED (S 15B 2N 0T)

DATA: U3411 #192

CALI: U3411 #2

BASE M/Z: 43

RIC: 12368.



100066

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: VERSAR INC.

Contract: _____

7

Lab Code: VERSAR Case No.: R3-7

SAS No.: _____ SDG No.: 1 _____

Matrix: (soil/water) SOIL

Lab Sample ID: 16431 _____

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: U3415 _____

Level: (low/med) LOW

Date Received: 04/19/90

% Moisture: not dec. 4

Date Analyzed: 04/27/90

Column: (pack/cap) PACK

Dilution Factor: 1.0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

Q

74-87-3-----Chloromethane	10	U
74-83-9-----Bromomethane	10	U
75-01-4-----Vinyl chloride	10	U
75-00-3-----Chloroethane	10	U
75-09-2-----Methylene chloride	5	U
67-64-1-----Acetone	10	U
75-15-0-----Carbon disulfide	5	U
75-35-4-----1,1-Dichloroethene	5	U
75-34-3-----1,1-Dichloroethane	5	U
540-59-0-----1,2-Dichloroethene (total)	5	U
67-66-3-----Chloroform	5	U
107-06-2-----1,2-Dichloroethane	5	U
78-93-3-----2-Butanone	10	U
71-55-6-----1,1,1-Trichloroethane	5	U
56-23-5-----Carbon tetrachloride	5	U
108-05-4-----Vinyl acetate	10	U
75-27-4-----Bromodichloromethane	5	U
78-87-5-----1,2-Dichloropropane	5	U
10061-01-5-----cis-1,3-Dichloropropene	5	U
79-01-6-----Trichloroethene	5	U
124-48-1-----Dibromochloromethane	5	U
79-00-5-----1,1,2-Trichloroethane	5	U
71-43-2-----Benzene	5	U
10061-02-6-----Trans-1,3-dichloropropene	5	U
75-25-2-----Bromoform	5	U
108-10-1-----4-Methyl-2-pentanone	10	U
591-78-6-----2-Hexanone	10	U
127-18-4-----Tetrachloroethene	5	U
79-34-5-----1,1,2,2-Tetrachloroethane	5	U
108-88-3-----Toluene	5	U
108-90-7-----Chlorobenzene	5	U
100-41-4-----Ethylbenzene	5	U
100-42-5-----Styrene	5	U
1330-20-7-----Total xylenes	5	U

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ORIGINAL
7 (Red)

Lab Name: VERSAR INC.

Contract: _____

Lab Code: VERSAR Case No.: R3-7 SAS No.: _____ SDG No.: 1 _____

Matrix: (soil/water) SOIL Lab Sample ID: 16431 _____

Sample wt/vol: 5.0 (g/mL) G Lab File ID: U3415 _____

Level: (low/med) LOW Date Received: 04/19/90

% Moisture: not dec. 4 Date Analyzed: 04/27/90

Column (pack/cap) PACK Dilution Factor: 1.0 _____

Number TICs found: 0

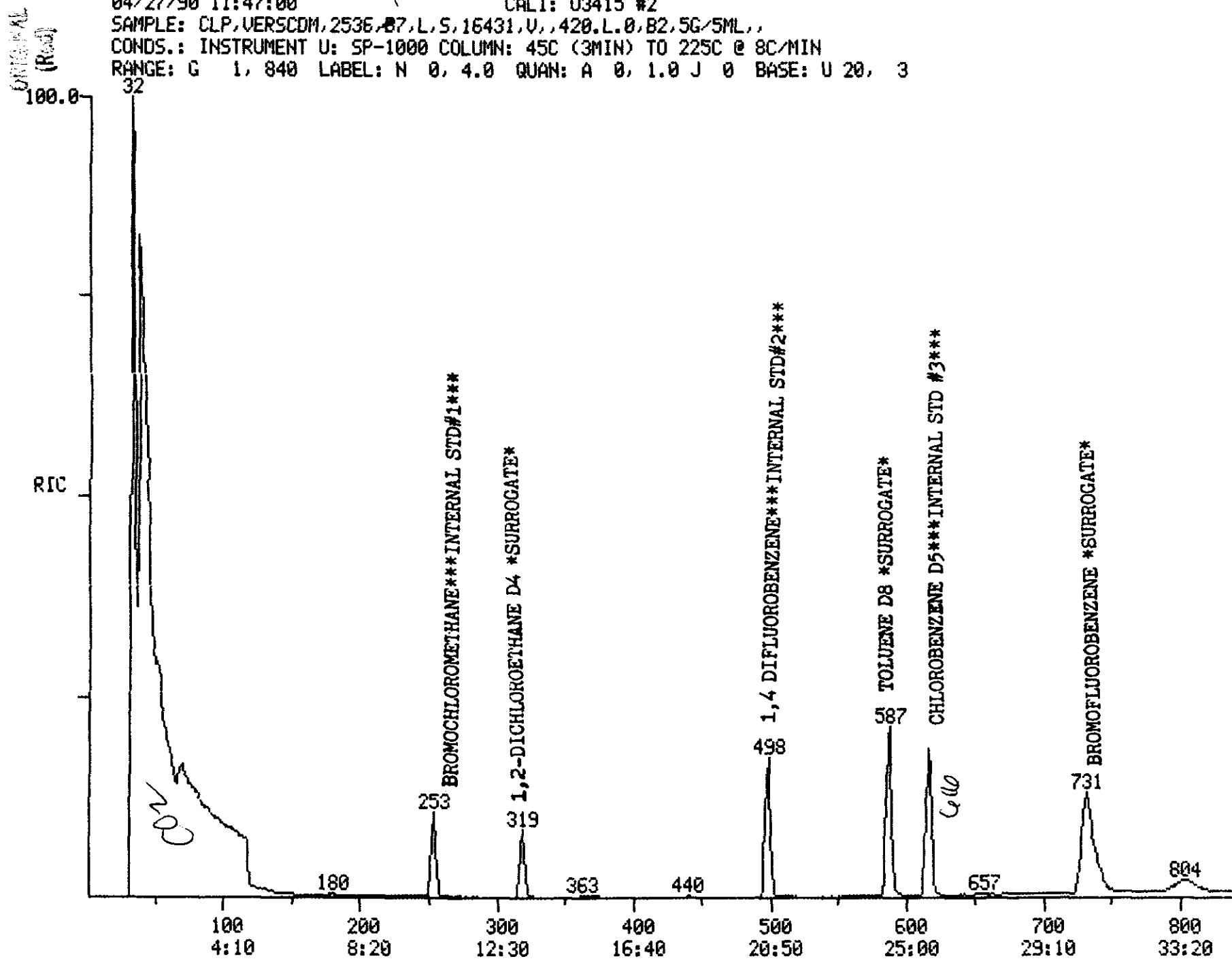
CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

RIC
04/27/90 11:47:00
SAMPLE: CLP, UERSO
COND.: INSTRUMENT
RANGE: G 1, 840
33

DATA: U3415 #1 SCANS 1 TO 840
 CALI: U3415 #2
 31.U., 420.L.0, B2.5G/5ML.,
 IN: 45C (3MIN) TU 225C @ 8C/MIN
 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

69000
678912.



Data: U3415.TI

04/27/90 11:47:00

Sample: CLP, VERSCDM, 2536, 87, L, S, 16431, V, , 420. L. O, B2, 5G/5ML, ,

Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

Formula: -

Instrument: U

Weight: 0.011

Submitted by: VERSAR

Analyst: MAT

Acct. No.: 420.1.0

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)

4.2 %

Resp. fac. from Library Entry

No	Name
1	C101 BROMOCHLOROMETHANE **IS#1**
2	C010 CHLOROMETHANE
3	C015 BROMOMETHANE
4	C020 VINYL CHLORIDE
5	C025 CHLOROETHANE
6	C030 METHYLENE CHLORIDE
7	C035 ACETONE
8	C040 CARBON DISULFIDE
9	C045 1,1-DICHLOROETHENE
10	C050 1,1-DICHLOROETHANE
11	C053 1,2-DICHLOROETHENE (TOTAL)
12	C060 CHLOROFORM
13	C065 1,2-DICHLOROETHANE
14	CS15 1,2-DICHLOROETHANE-D4 **SS#1**
15	C043 TRICHLOROFLUOROMETHANE
16	CI10 1,4-DIFLUOROBENZENE **IS#2**
17	C110 2-BUTANONE
18	C125 VINYL ACETATE
19	C120 CARBON TETRACHLORIDE
20	C130 BROMODICHLOROMETHANE
21	C140 1,2-DICHLOROPROPANE
22	C145 CIS-1,3-DICHLOROPROPENE
23	C150 TRICHLOROETHENE
24	C165 BENZENE
25	C155 DIBROMOCHLOROMETHANE
26	C160 1,1,2-TRICHLOROETHANE
27	C170 TRANS-1,3-DICHLOROPROPENE
28	C180 BROMOFORM
29	C115 1,1,1-TRICHLOROETHANE
30	CS05 TOLUENE-D8 **SS#2**
31	CI20 CHLOROBENZENE-D5 **IS#3**
32	C205 4-METHYL-2-PENTANONE
33	C210 2-HEXANONE
34	C220 TETRACHLOROETHENE
35	C225 1,1,2,2-TETRACHLOROETHANE
36	C230 TOLUENE
37	C235 CHLOROBENZENE
38	C240 ETHYLBENZENE
39	C245 STYRENE
40	C250 ORTHO & PARA XYLENE
41	C251 META-XYLENE
42	CS10 BROMOFLUOROBENZENE **SS#3**

INT STD COMPLIANT

SURROG. COMPLIANT

M.A. Templeton

4/27/90

100070

u3415

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	128	253	10:32	1	1.000	A BB	49516.	50.000 UG/L*	16.72
2	NOT FOUND								ORIGINAL
3	NOT FOUND								(Red)
4	NOT FOUND								
5	NOT FOUND								
6	84	179	7:27	1	0.708	A BB	2305.	1.085 UG/L	0.61
7	43	193	8:02	1	0.763	A BB	480.	0.981 UG/L	0.33
8	NOT FOUND								
9	NOT FOUND								
10	NOT FOUND								
11	NOT FOUND								
12	NOT FOUND								
13	NOT FOUND								
14	65	318	13:15	1	1.257	A BB	86229.	43.355 UG/L%	14.50 87%
15	NOT FOUND								
16	114	498	20:45	16	1.000	A BB	216099.	50.000 UG/L*	16.72
17	NOT FOUND								
18	NOT FOUND								
19	NOT FOUND								
20	NOT FOUND								
21	NOT FOUND								
22	NOT FOUND								
23	NOT FOUND								
24	NOT FOUND								
25	NOT FOUND								
26	NOT FOUND								
27	NOT FOUND								
28	NOT FOUND								
29	NOT FOUND								
30	98	587	24:27	31	0.953	A BB	231311.	55.328 UG/L%	18.50 11%
31	117	616	25:40	31	1.000	A BB	181858.	50.000 UG/L*	16.72
32	NOT FOUND								
33	NOT FOUND								
34	NOT FOUND								
35	NOT FOUND								
36	92	591	24:37	31	0.959	A BB	545.	0.235 UG/L	0.08
37	NOT FOUND								
38	NOT FOUND								
39	NOT FOUND								
40	NOT FOUND								
41	NOT FOUND								
42	95	731	30:27	31	1.187	A BB	129587.	47.302 UG/L%	15.82 95%

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
1	10:30	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
2	2:10		0.206						
3	3:20		0.317						
4	4:12		0.401						
5	5:15		0.500						
6	7:25	1.01	0.706	1.00	1.83	50.00	0.047	1.268	0.04
7	8:00	1.01	0.762	1.00	0.98	50.00	0.010	0.494	0.02
8	9:00		0.857						
9	10:05		0.960						
10	11:22		1.083						
11	12:02		1.147						
12	12:37		1.202						

100071

u3415
ORIGINAL
(Rev)

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
13	13:20		1. 270						
14	13:15	1. 00	1. 262	1. 00	43. 35	50. 00	1. 741	2. 008	0. 87
15	9:27		0. 901						
16	20:42	1. 00	1. 000	1. 00	50. 00	50. 00	1. 000	1. 000	1. 00
17	13:15		1. 262						
18	15:07		0. 730						
19	15:05		0. 728						
20	15:30		0. 748						
21	16:55		0. 817						
22	17:07		0. 827						
23	17:40		0. 853						
24	18:12		0. 879						
25	18:17		0. 883						
26	18:25		0. 889						
27	18:22		0. 887						
28	20:57		1. 012						
29	14:42		0. 710						
30	24:27	1. 00	0. 953	1. 00	55. 33	50. 00	1. 272	1. 149	1. 11
31	25:40	1. 00	1. 000	1. 00	50. 00	50. 00	1. 000	1. 000	1. 00
32	21:30		0. 838						
33	23:02		0. 898						
34	23:17		0. 907						
35	23:17		0. 907						
36	24:37	1. 00	0. 959	1. 00	0. 23	50. 00	0. 003	0. 638	0. 00
37	25:47		1. 005						
38	27:57		1. 089						
39	32:25		1. 263						
40	33:47		1. 317						
41	32:42		1. 274						
42	30:45	0. 99	1. 198	0. 99	47. 30	50. 00	0. 713	0. 753	0. 95

100072

OKC(1-83)
(Rev A)

Versar^{inc.}

IV. STANDARDS DATA

100073

6A
VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: VERSAR INC.

Contract: _____

La Code: VERSAR Case No.: R3-7

SAS No.: _____ SDG No.: 1

Instrument ID: U

Calibration Date(s): 04/23/90 04/23/90

Matrix: (soil/water) SOIL Level: (low/med) LOW Column: (pack/cap) PACK

Min RRF for SPCC(#) = 0.300 (0.250 for Bromoform) Max %RSD for CCC(*) = 30.0%

LAB FILE ID:	RRF20 = <u>U3331</u>	RRF50 = <u>U3327</u>
RRF100= <u>U3330</u>	RRF150= <u>U3329</u>	RRF200= <u>U3328</u>

COMPOUND	RRF20	RRF50	RRF100	RRF150	RRF200	<u>RRF</u>	% RSD
Chloromethane	# 1.080	1.389	1.256	1.475	1.362	1.312	11.6#
Propanomethane	0.840	0.860	0.833	0.616	0.502	0.730	22.2
Methyl chloride	* 1.027	1.115	1.140	1.009	0.973	1.053	6.8*
Chloroethane	0.549	0.601	0.608	0.577	0.606	0.588	4.3
Methylene chloride	1.147	1.277	1.250	1.317	1.299	1.258	5.3
Acetone	0.271	0.442	0.287	0.340	0.352	0.338	19.9
Carbon disulfide	0.937	1.267	1.120	1.240	1.304	1.174	12.7
1,1-Dichloroethene	* 0.977	1.052	1.160	1.222	1.212	1.125	9.5*
1,1-Dichloroethane	# 1.928	2.229	2.253	2.414	2.444	2.254	9.1#
1,2-Dichloroethene (total)	1.090	1.130	1.357	1.343	1.327	1.249	10.3
Chloroform	* 2.261	2.495	2.506	2.605	2.581	2.490	5.5*
1,1-Dichloroethane	1.402	1.784	1.766	1.851	1.898	1.740	11.3
2-Butanone	0.172	0.217	0.190	0.216	0.219	0.203	10.3
1,1,1-Trichloroethane	0.419	0.404	0.499	0.483	0.480	0.457	9.3
Carbon tetrachloride	0.399	0.378	0.448	0.441	0.430	0.419	7.1
Vinyl acetate	0.389	0.558	0.544	0.614	0.686	0.558	19.7
Bromodichloromethane	0.467	0.511	0.568	0.550	0.550	0.529	7.7
1,2-Dichloropropane	* 0.302	0.333	0.345	0.367	0.367	0.343	7.9*
1,1,3-Dichloropropene	0.556	0.600	0.680	0.658	0.653	0.629	8.0
Trichloroethene	0.427	0.412	0.447	0.450	0.430	0.433	3.6
Dibromochloromethane	0.473	0.464	0.481	0.455	0.424	0.459	4.8
1,1,2-Trichloroethane	0.309	0.315	0.319	0.322	0.314	0.316	1.6
Benzene	0.815	0.827	0.936	0.903	0.886	0.873	5.9
Trans-1,3-dichloropropene	0.294	0.320	0.362	0.346	0.346	0.334	8.0
Bromoform	# 0.427	0.433	0.508	0.476	0.454	0.460	7.2#
4-Methyl-2-pentanone	0.452	0.652	0.550	0.665	0.723	0.608	17.7
2-Hexanone	0.399	0.540	0.465	0.565	0.567	0.507	14.4
Tetrachloroethene	0.446	0.399	0.431	0.417	0.386	0.416	5.8
1,1,2,2-Tetrachloroethane	# 0.645	0.683	0.704	0.699	0.675	0.681	3.4#
Toluene	* 0.620	0.621	0.709	0.703	0.677	0.666	6.5*
Chlorobenzene	# 0.870	0.876	0.901	0.922	0.892	0.892	2.3#
Ethylbenzene	* 0.406	0.396	0.474	0.470	0.452	0.440	8.3*
Styrene	0.776	0.801	0.796	0.863	0.903	0.828	6.4
Total xylenes	0.438	0.448	0.452	0.497	0.524	0.472	7.8
m-Xylene	0.501	0.519	0.523	0.566	0.599	0.542	7.4
Truene-d8	0.960	1.213	1.076	1.082	1.019	1.070	8.8
Bromofluorobenzene	0.694	0.832	0.755	0.767	0.731	0.756	6.7
1,2-Dichloroethane-d4	1.279	1.790	1.546	1.632	1.648	1.579	12.0

7A
VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: VERSAR INC. Contract: _____
 La Code: VERSAR Case No.: R3-7 SAS No.: _____ SDG No.: 1
 Instrument ID: U Calibration date: 04/26/90 Time: 830
 Lab File ID: U3386 Init. Calib. Date(s): 04/23/90 04/23/90
 Matrix: (soil/water) SOIL Level: (low/med) LOW Column: (pack/cap) PACK
 Min RRF50 for SPCC(#) = 0.300 (0.250 for Bromoform) Max %D for CCC(*) = 25.0%

COMPOUND	RRF	RRF50	%D
Chloromethane	# 1.312	1.348	-2.7 #
Bromomethane	0.730	0.855	-17.1
Vinyl chloride	* 1.053	1.019	3.2 *
Chloroethane	0.588	0.682	-16.0
Methylene chloride	1.258	1.372	-9.1
Acetone	0.338	0.727	-99.9
Carbon disulfide	1.174	1.045	11.0
1,1-Dichloroethene	* 1.125	1.152	-2.4 *
1,1-Dichloroethane	# 2.254	2.837	-25.9 #
1,2-Dichloroethene (total)	1.249	1.354	-8.4
Chloroform	* 2.490	2.959	-18.8 *
1,2-Dichloroethane	1.740	2.513	-44.4
2-Butanone	0.203	0.236	-16.3
1,1,1-Trichloroethane	0.457	0.493	-7.9
Carbon tetrachloride	0.419	0.428	-2.1
Vinyl acetate	0.558	0.791	-41.8
Bromodichloromethane	0.529	0.580	-9.6
1,2-Dichloropropane	* 0.343	0.392	-14.3 *
cis-1,3-Dichloropropene	0.629	0.679	-7.9
Trichloroethene	0.433	0.433	0.0
Dibromochloromethane	0.459	0.486	-5.9
1,1,2-Trichloroethane	0.316	0.352	-11.4
Benzene	0.873	0.889	-1.8
Trans-1,3-dichloropropene	0.334	0.366	-9.6
Bromoform	# 0.460	0.446	3.0 #
4-Methyl-2-pentanone	0.608	0.922	-51.6
2-Hexanone	0.507	0.759	-49.7
Tetrachloroethene	0.416	0.417	-0.2
1,1,2,2-Tetrachloroethane	# 0.681	0.713	-4.7 #
Toluene	* 0.666	0.686	-3.0 *
Chlorobenzene	# 0.892	0.915	-2.6 #
Ethylbenzene	* 0.440	0.439	0.2 *
Styrene	0.828	0.862	-4.1
Total xylenes	0.472	0.477	-1.1
m-Xylene	0.542	0.559	-3.1
Toluene-d8	1.070	1.202	-12.3
Bromofluorobenzene	0.756	0.869	-15.0
1,2-Dichloroethane-d4	1.579	2.288	-44.9

7A
VOLATILE CONTINUING CALIBRATION CHECK

(RGA)

Lab Name: VERSAR INC.

Contract: _____

La. Code: VERSAR Case No.: R3-7 SAS No.: _____ SDG No.: 1 _____

Instrument ID: U Calibration date: 04/27/90 Time: 821 _____

Lab File ID: U3411 Init. Calib. Date(s): 04/23/90 04/23/90

Matrix: (soil/water) SOIL Level: (low/med) LOW Column: (pack/cap) PACK

Min RRF50 for SPCC(#) = 0.300 (0.250 for Bromoform) Max %D for CCC(*) = 25.0%

COMPOUND	RRF	RRF50	%D
Chloromethane	# 1.312	1.321	-0.7 #
Bromomethane	0.730	1.030	-41.1
Vinyl chloride	* 1.053	1.057	-0.4 *
Chloroethane	0.588	0.776	-32.0
Methylene chloride	1.258	1.268	-0.8
Acetone	0.338	0.494	-46.2
Carbon disulfide	1.174	0.784	33.2
1,1-Dichloroethene	* 1.125	0.992	11.8 *
1,1-Dichloroethane	# 2.254	2.426	-7.6 #
1,2-Dichloroethene (total)	1.249	1.193	4.5
Chloroform	* 2.490	2.598	-4.3 *
1,2-Dichloroethane	1.740	2.119	-21.8
2-Butanone	0.203	0.197	3.0
1,1,1-Trichloroethane	0.457	0.496	-8.5
Carbon tetrachloride	0.419	0.436	-4.1
Vinyl acetate	0.558	0.619	-10.9
Bromodichloromethane	0.529	0.585	-10.6
1,2-Dichloropropane	* 0.343	0.382	-11.4 *
cis-1,3-Dichloropropene	0.629	0.675	-7.3
Trichloroethene	0.433	0.466	-7.6
Dibromochloromethane	0.459	0.527	-14.8
1,1,2-Trichloroethane	0.316	0.342	-8.2
Benzene	0.873	0.901	-3.2
Trans-1,3-dichloropropene	0.334	0.361	-8.1
Bromoform	# 0.460	0.493	-7.2 #
4-Methyl-2-pentanone	0.608	0.725	-19.2
2-Hexanone	0.507	0.579	-14.2
Tetrachloroethene	0.416	0.481	-15.6
1,1,2,2-Tetrachloroethane	# 0.681	0.703	-3.2 #
Toluene	* 0.666	0.638	4.2 *
Chlorobenzene	# 0.892	0.892	0.0 #
Ethylbenzene	* 0.440	0.411	6.6 *
Styrene	0.828	0.770	7.0
Total xylenes	0.472	0.438	7.2
m-Xylene	0.542	0.505	6.8
Toluene-d8	1.070	1.149	-7.4
Bromofluorobenzene	0.756	0.753	0.4
1,2-Dichloroethane-d4	1.579	2.008	-27.2

RIC
04/23/90 19:35:00

DATA: U3331 #1
CALI: U3331 #2

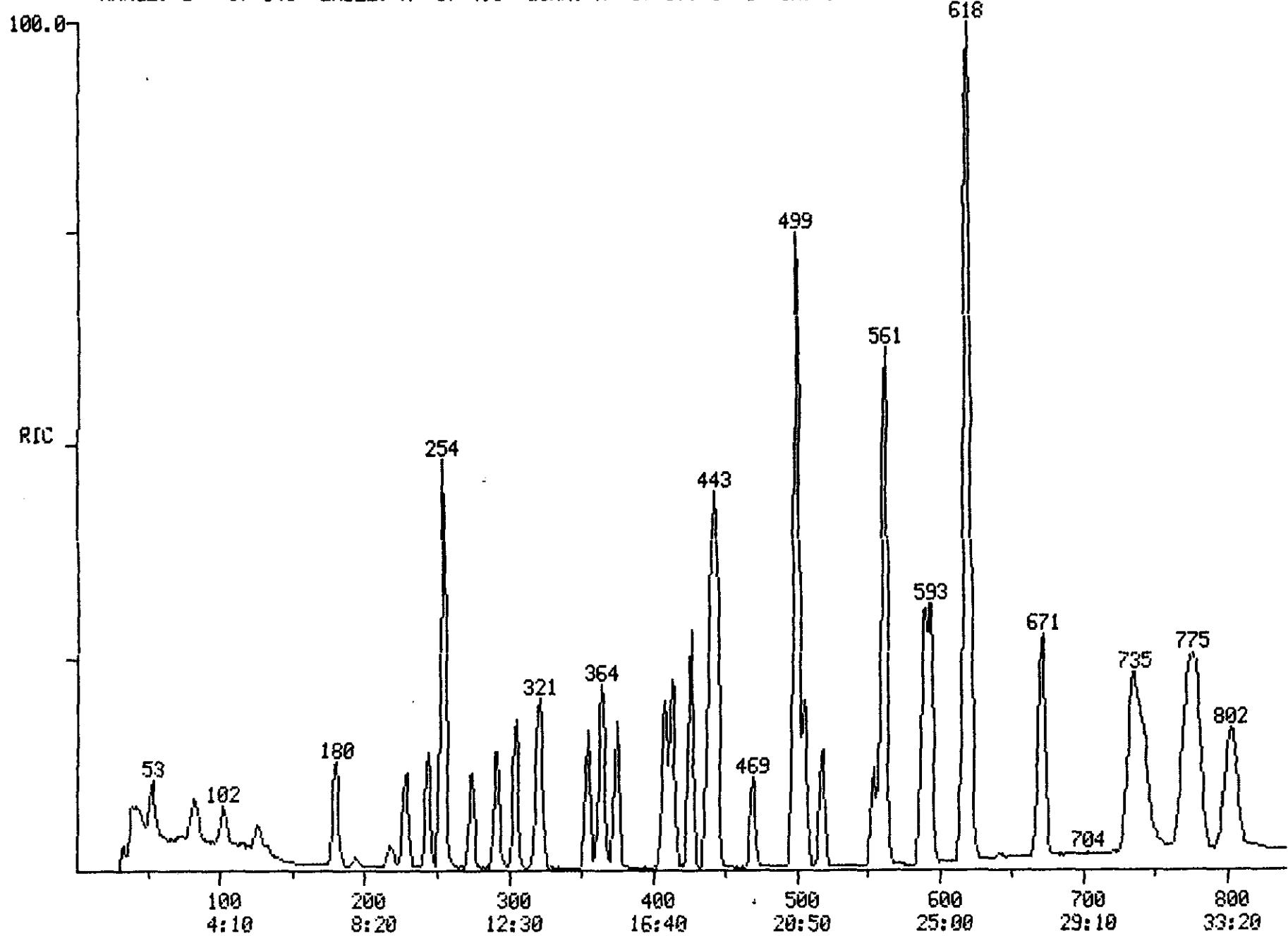
SCANS 1 TO 840

SAMPLE: CLP,,,VSTD20,L,S,22526,V,IC-020,,,5ML,

COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

RANGE: G 1, 840 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

155136.



Quantitation Report File: U3331

ORIGINAL

(R: 0)

Data: U3331.TI

04/23/90 19:35:00

Sample: CLP,,, VSTD20, L, S, 22526, V, IC-020,,, 5ML,

Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

Formula: - Instrument: U Weight: 0.012

Submitted by: VERSAR Analyst: DD

Acct. No.: -

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)

Resp. fac. from Library Entry

No	Name
1	C101 BROMOCHLOROMETHANE **IS#1**
2	C010 CHLOROMETHANE
3	C015 BROMOMETHANE
4	C020 VINYL CHLORIDE
5	C025 CHLOROETHANE
6	C030 METHYLENE CHLORIDE
7	C035 ACETONE
8	C040 CARBON DISULFIDE
9	C045 1,1-DICHLOROETHENE
10	C050 1,1-DICHLOROETHANE
11	C053 1,2-DICHLOROETHENE (TOTAL)
12	C060 CHLOROFORM
13	C065 1,2-DICHLOROETHANE
14	CS15 1,2-DICHLOROETHANE-D4 **SS#1**
15	C043 TRICHLOROFLUOROMETHANE
16	CI10 1,4-DIFLUOROBENZENE **IS#2**
17	C110 2-BUTANONE
18	C125 VINYL ACETATE
19	C120 CARBON TETRACHLORIDE
20	C130 BROMODICHLOROMETHANE
21	C140 1,2-DICHLOROPROPANE
22	C145 CIS-1,3-DICHLOROPROPENE
23	C150 TRICHLOROETHENE
24	C165 BENZENE
25	C155 DIBROMOCHLOROMETHANE
26	C160 1,1,2-TRICHLOROETHANE
27	C170 TRANS-1,3-DICHLOROPROPENE
28	C180 BROMOFORM
29	C115 1,1,1-TRICHLOROETHANE
30	CS05 TOLUENE-DB **SS#2**
31	CI20 CHLOROBENZENE-D5 **IS#3**
32	C205 4-METHYL-2-PENTANONE
33	C210 2-HEXANONE
34	C220 TETRACHLOROETHENE
35	C225 1,1,2,2-TETRACHLOROETHANE
36	C230 TOLUENE
37	C235 CHLOROBENZENE
38	C240 ETHYLBENZENE
39	C245 STYRENE
40	C250 ORTHO & PARA XYLENE
41	C251 META-XYLENE
42	CS10 BROMOFLUOROBENZENE **SS#3**

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U 33 ORIGINAL
(Part)

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	128	254	10:35	1	1.000	A BB	50907.	50.000 UG/L*	5.38
2	50	53	2:12	1	0.209	A BB	22001.	20.000 UG/L	2.15
3	94	82	3:25	1	0.323	A BB	17104.	20.000 UG/L	2.15
4	62	102	4:15	1	0.402	A BB	20913.	20.000 UG/L	2.15
5	64	126	5:15	1	0.496	A BB	11174.	20.000 UG/L	2.15
6	84	179	7:27	1	0.705	A BB	23355.	20.000 UG/L	2.15
7	43	194	8:05	1	0.764	A BB	5517.	20.000 UG/L	2.15
8	76	218	9:05	1	0.858	A BB	19071.	20.000 UG/L	2.15
9	96	244	10:10	1	0.961	A BB	19888.	20.000 UG/L	2.15
10	63	275	11:27	1	1.083	A BB	39265.	20.000 UG/L	2.15
11	96	291	12:07	1	1.146	A BB	22195.	20.000 UG/L	2.15
12	83	305	12:42	1	1.201	A BB	46049.	20.000 UG/L	2.15
13	62	322	13:25	1	1.268	A BB	28541.	20.000 UG/L%	2.15
14	65	320	13:20	1	1.260	A BB	26036.	20.000 UG/L	2.15
15	101	228	9:30	1	0.898	A BB	38813.	20.000 UG/L	2.15
16	114	499	20:47	16	1.000	A BB	225674.	50.000 UG/L*	5.38
17	72	320	13:20	1	1.260	A BB	3498.	20.000 UG/L	2.15
18	43	366	15:15	16	0.733	A BB	35087.	20.000 UG/L	2.15
19	117	364	15:10	16	0.729	A VB	36013.	20.000 UG/L	2.15
20	83	375	15:37	16	0.752	A BB	42132.	20.000 UG/L	2.15
21	63	408	17:00	16	0.818	A BB	27305.	20.000 UG/L	2.15
22	75	413	17:12	16	0.828	A BB	50165.	20.000 UG/L	2.15
23	130	426	17:45	16	0.854	A BB	38565.	20.000 UG/L	2.15
24	78	439	18:17	16	0.880	A BB	73586.	20.000 UG/L	2.15
25	129	441	18:22	16	0.884	A BB	42725.	20.000 UG/L	2.15
26	97	444	18:30	16	0.890	A BB	27937.	20.000 UG/L	2.15
27	75	444	18:30	16	0.890	A BB	26535.	20.000 UG/L	2.15
28	173	505	21:02	16	1.012	A BB	38545.	20.000 UG/L	2.15
29	97	355	14:47	16	0.711	A BB	37856.	20.000 UG/L	2.15
30	98	588	24:30	31	0.953	A BB	80925.	20.000 UG/L%	2.15
31	117	617	25:42	31	1.000	A BB	210646.	50.000 UG/L*	5.38
32	43	518	21:35	31	0.840	A BB	38088.	20.000 UG/L	2.15
33	43	554	23:05	31	0.898	A BB	33600.	20.000 UG/L	2.15
34	164	561	23:22	31	0.909	A BB	37555.	20.000 UG/L	2.15
35	83	561	23:22	31	0.909	A BB	54357.	20.000 UG/L	2.15
36	92	593	24:42	31	0.961	A BB	52259.	20.000 UG/L	2.15
37	112	620	25:50	31	1.005	A BB	73346.	20.000 UG/L	2.15
38	106	670	27:55	31	1.086	A BB	34177.	20.000 UG/L	2.15
39	104	772	32:10	31	1.251	A BB	65388.	20.000 UG/L	2.15
40	106	802	33:25	31	1.300	A BB	36925.	20.000 UG/L	2.15
41	106	777	32:22	31	1.259	A BB	42193.	20.000 UG/L	2.15
42	95	734	30:35	31	1.190	A BB	58456.	20.000 UG/L%	2.15

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
1	10:35	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
2	2:12	1.00	0.209	1.00	20.00	20.00	1.080	1.080	1.00
3	3:25	1.00	0.323	1.00	20.00	20.00	0.840	0.840	1.00
4	4:15	1.00	0.402	1.00	20.00	20.00	1.027	1.027	1.00
5	5:15	1.00	0.496	1.00	20.00	20.00	0.549	0.549	1.00
6	7:27	1.00	0.705	1.00	20.00	20.00	1.147	1.147	1.00
7	8:05	1.00	0.764	1.00	20.00	20.00	0.271	0.271	1.00
8	9:05	1.00	0.858	1.00	20.00	20.00	0.937	0.937	1.00
9	10:10	1.00	0.961	1.00	20.00	20.00	0.977	0.977	1.00
10	11:27	1.00	1.083	1.00	20.00	20.00	1.090	1.090	1.00
11	12:07	1.00	1.146	1.00	20.00	20.00	2.261	2.261	1.00
12	12:42	1.00	1.201	1.00	20.00	20.00			

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No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R.Fac	R.Fac(L)	Ratio
(Ref)									
13	13:25	1.00	1.268	1.00	20.00	20.00	1.402	1.402	1.00
14	13:20	1.00	1.260	1.00	20.00	20.00	1.279	1.279	1.00
15	9:30	1.00	0.898	1.00	20.00	20.00	1.906	1.906	1.00
16	20:47	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
17	13:20	1.00	1.260	1.00	20.00	20.00	0.172	0.172	1.00
18	15:15	1.00	0.733	1.00	20.00	20.00	0.389	0.389	1.00
19	15:10	1.00	0.729	1.00	20.00	20.00	0.399	0.399	1.00
20	15:37	1.00	0.752	1.00	20.00	20.00	0.467	0.467	1.00
21	17:00	1.00	0.818	1.00	20.00	20.00	0.302	0.302	1.00
22	17:12	1.00	0.828	1.00	20.00	20.00	0.556	0.556	1.00
23	17:45	1.00	0.854	1.00	20.00	20.00	0.427	0.427	1.00
24	18:17	1.00	0.880	1.00	20.00	20.00	0.815	0.815	1.00
25	18:22	1.00	0.884	1.00	20.00	20.00	0.473	0.473	1.00
26	18:30	1.00	0.890	1.00	20.00	20.00	0.309	0.309	1.00
27	18:30	1.00	0.890	1.00	20.00	20.00	0.294	0.294	1.00
28	21:02	1.00	1.012	1.00	20.00	20.00	0.427	0.427	1.00
29	14:47	1.00	0.711	1.00	20.00	20.00	0.419	0.419	1.00
30	24:30	1.00	0.953	1.00	20.00	20.00	0.960	0.960	1.00
31	25:42	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
32	21:35	1.00	0.840	1.00	20.00	20.00	0.452	0.452	1.00
33	23:05	1.00	0.898	1.00	20.00	20.00	0.399	0.399	1.00
34	23:22	1.00	0.909	1.00	20.00	20.00	0.446	0.446	1.00
35	23:22	1.00	0.909	1.00	20.00	20.00	0.645	0.645	1.00
36	24:42	1.00	0.961	1.00	20.00	20.00	0.620	0.620	1.00
37	25:50	1.00	1.005	1.00	20.00	20.00	0.870	0.870	1.00
38	27:55	1.00	1.086	1.00	20.00	20.00	0.406	0.406	1.00
39	32:10	1.00	1.251	1.00	20.00	20.00	0.776	0.776	1.00
40	33:25	1.00	1.300	1.00	20.00	20.00	0.438	0.438	1.00
41	32:22	1.00	1.259	1.00	20.00	20.00	0.504	0.504	1.00
42	30:35	1.00	1.190	1.00	20.00	20.00	0.694	0.694	1.00

REF ID: A

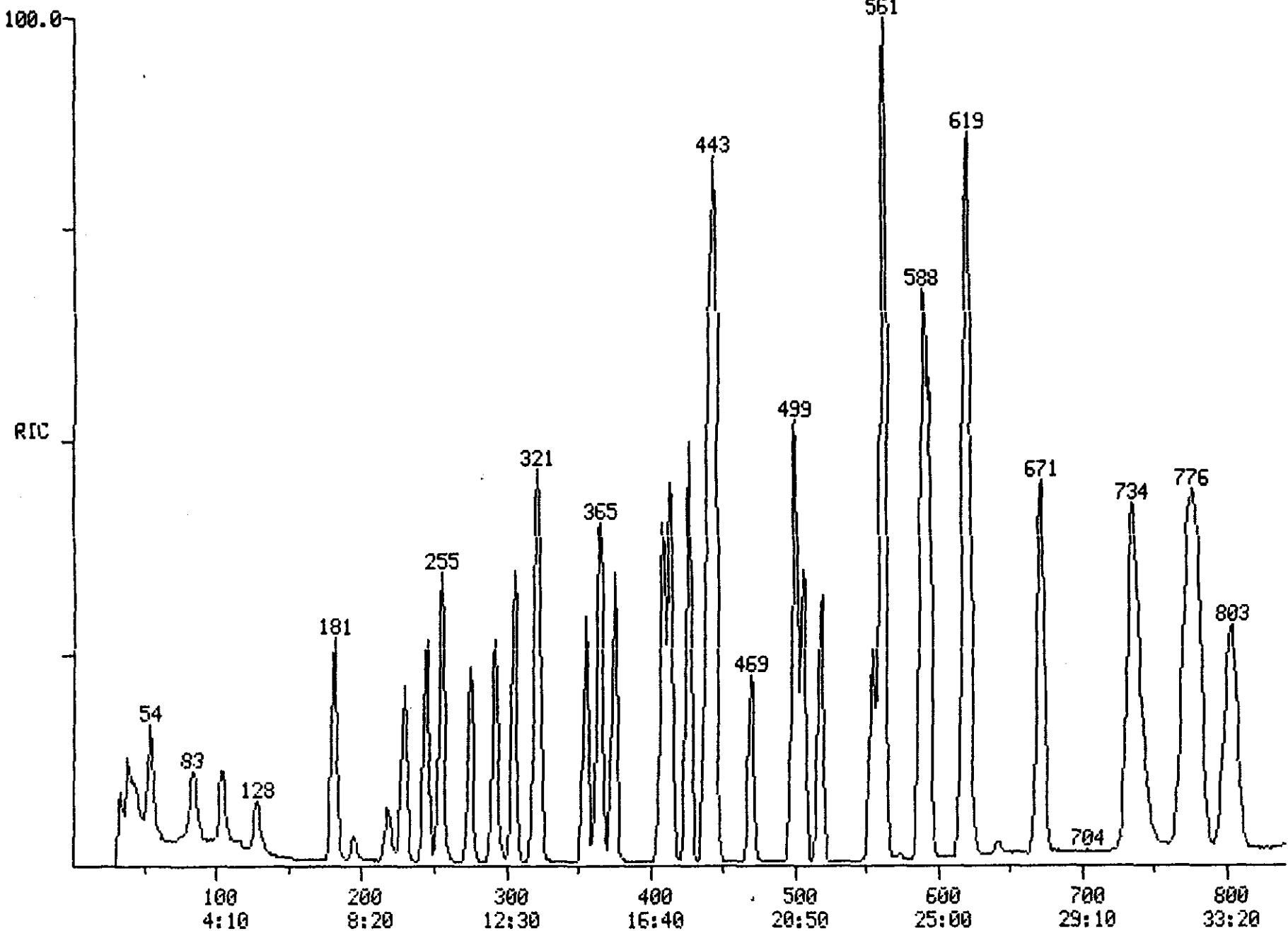
ORIGINAL
(Ref)

RIC
04/23/90 16:23:00
SAMPLE: CLP.,,VSTD50,L,S,22522,U,IC-050,,,5ML,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
RANGE: G 1, 840 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: U3327 #1
CALI: U3327 #2

SCANS 1 TO 840

24115
N00081



Data: U3327.TI

04/23/90 16:23:00

Sample: CLP,,,VSTD50,L,S,22522,V,IC-050,,,5ML,

Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

Formula: - Instrument: U Weight: 0.012

Submitted by: VERSAR Analyst: DD Acct. No.: -

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)

Resp. fac. from Library Entry

No	Name
1	C101 BROMOCHLOROMETHANE **IS#1**
2	C010 CHLOROMETHANE
3	C015 BROMOMETHANE
4	C020 VINYL CHLORIDE
5	C025 CHLOROETHANE
6	C030 METHYLENE CHLORIDE
7	C035 ACETONE
8	C040 CARBON DISULFIDE
9	C045 1,1-DICHLOROETHENE
10	C050 1,1-DICHLOROETHANE
11	C053 1,2-DICHLOROETHENE (TOTAL)
12	C060 CHLOROFORM
13	C065 1,2-DICHLOROETHANE
14	CS15 1,2-DICHLOROETHANE-D4 **SS#1**
15	C043 TRICHLOROFUOROMETHANE
16	C110 1,4-DIFLUOROBENZENE **IS#2**
17	C110 2-BUTANONE
18	C125 VINYL ACETATE
19	C120 CARBON TETRACHLORIDE
20	C130 BROMODICHLOROMETHANE
21	C140 1,2-DICHLOROPROPANE
22	C145 CIS-1,3-DICHLOROPROPENE
23	C150 TRICHLOROETHENE
24	C165 BENZENE
25	C155 DIBROMOCHLOROMETHANE
26	C160 1,1,2-TRICHLOROETHANE
27	C170 TRANS-1,3-DICHLOROPROPENE
28	C180 BROMOFORM
29	C115 1,1,1-TRICHLOROETHANE
30	CS05 TOLUENE-D8 **SS#2**
31	C120 CHLOROBENZENE-D5 **IS#3**
32	C205 4-METHYL-2-PENTANONE
33	C210 2-HEXANONE
34	C220 TETRACHLOROETHENE
35	C225 1,1,2,2-TETRACHLOROETHANE
36	C230 TOLUENE
37	C235 CHLOROBENZENE
38	C240 ETHYLBENZENE
39	C245 STYRENE
40	C250 ORTHO & PARA XYLENE
41	C251 META-XYLENE
42	CS10 BROMOFLUOROBENZENE **SS#3**

(R.E.)

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	UG/L*	% Tot
1	128	255	10:37	1	1.000	A BB	52222.	50.000	UG/L*	2.38
2	50	54	2:15	1	0.212	A BB	72540.	50.000	UG/L	2.38
3	94	83	3:27	1	0.325	A BB	44892.	50.000	UG/L	2.38
4	62	104	4:20	1	0.408	A BB	58220.	50.000	UG/L	2.38
5	64	128	5:20	1	0.502	A BB	31368.	50.000	UG/L	2.38
6	84	181	7:32	1	0.710	A BB	66686.	50.000	UG/L	2.38
7	43	195	8:07	1	0.765	A BB	23081.	50.000	UG/L	2.38
8	76	218	9:05	1	0.855	A BB	66183.	50.000	UG/L	2.38
9	96	245	10:12	1	0.961	A BB	54961.	50.000	UG/L	2.38
10	63	275	11:27	1	1.078	A BB	116419.	50.000	UG/L	2.38
11	96	292	12:10	1	1.145	A BB	59012.	50.000	UG/L	2.38
12	83	305	12:42	1	1.196	A BB	130308.	50.000	UG/L	2.38
13	62	322	13:25	1	1.263	A BB	93181.	50.000	UG/L	2.38
14	65	320	13:20	1	1.255	A BB	93473.	50.000	UG/L%	2.38
15	101	230	9:35	1	0.902	A BB	103525.	50.000	UG/L	2.38
16	114	499	20:47	16	1.000	A BB	236574.	50.000	UG/L*	2.38
17	72	321	13:22	1	1.259	A BB	11317.	50.000	UG/L	2.38
18	43	366	15:15	16	0.733	A BB	131925.	50.000	UG/L	2.38
19	117	364	15:10	16	0.729	A VB	89337.	50.000	UG/L	2.38
20	83	375	15:37	16	0.752	A BB	120808.	50.000	UG/L	2.38
21	63	408	17:00	16	0.818	A BB	78730.	50.000	UG/L	2.38
22	75	413	17:12	16	0.828	A BB	142018.	50.000	UG/L	2.38
23	130	426	17:45	16	0.854	A BB	97488.	50.000	UG/L	2.38
24	78	439	18:17	16	0.880	A BB	195594.	50.000	UG/L	2.38
25	129	441	18:22	16	0.884	A BB	109733.	50.000	UG/L	2.38
26	97	444	18:30	16	0.890	A BB	74531.	50.000	UG/L	2.38
27	75	444	18:30	16	0.890	A BB	75643.	50.000	UG/L	2.38
28	173	505	21:02	16	1.012	A BB	102512.	50.000	UG/L	2.38
29	97	355	14:47	16	0.711	A BB	95527.	50.000	UG/L	2.38
30	98	588	24:30	31	0.953	A BB	262709.	50.000	UG/L%	2.38
31	117	617	25:42	31	1.000	A BB	216519.	50.000	UG/L*	2.38
32	43	518	21:35	31	0.840	A BB	141121.	50.000	UG/L	2.38
33	43	554	23:05	31	0.898	A BB	116928.	50.000	UG/L	2.38
34	164	561	23:22	31	0.909	A BB	86319.	50.000	UG/L	2.38
35	83	561	23:22	31	0.909	A BB	147916.	50.000	UG/L	2.38
36	92	593	24:42	31	0.961	A BB	134473.	50.000	UG/L	2.38
37	112	620	25:50	31	1.005	A BB	189769.	50.000	UG/L	2.38
38	106	670	27:55	31	1.086	A BB	85664.	50.000	UG/L	2.38
39	104	771	32:07	31	1.250	A BB	173441.	50.000	UG/L	2.38
40	106	803	33:27	31	1.301	A BB	97014.	50.000	UG/L	2.38
41	106	778	32:25	31	1.261	A BB	112396.	50.000	UG/L	2.38
42	95	734	30:35	31	1.190	A BB	180092.	50.000	UG/L%	2.38

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
1	10:37	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
2	2:15	1.00	0.212	1.00	50.00	50.00	1.389	1.389	1.00
3	3:27	1.00	0.325	1.00	50.00	50.00	0.860	0.860	1.00
4	4:20	1.00	0.408	1.00	50.00	50.00	1.115	1.115	1.00
5	5:20	1.00	0.502	1.00	50.00	50.00	0.601	0.601	1.00
6	7:32	1.00	0.710	1.00	50.00	50.00	1.277	1.277	1.00
7	8:07	1.00	0.765	1.00	50.00	50.00	0.442	0.442	1.00
8	9:05	1.00	0.855	1.00	50.00	50.00	1.267	1.267	1.00
9	10:12	1.00	0.961	1.00	50.00	50.00	1.052	1.052	1.00
10	11:27	1.00	1.078	1.00	50.00	50.00	2.229	2.229	1.00
11	12:10	1.00	1.145	1.00	50.00	50.00	1.130	1.130	1.00
12	12:42	1.00	1.196	1.00	50.00	50.00	2.495	2.495	1.00

100083

00000000000000000000000000000000

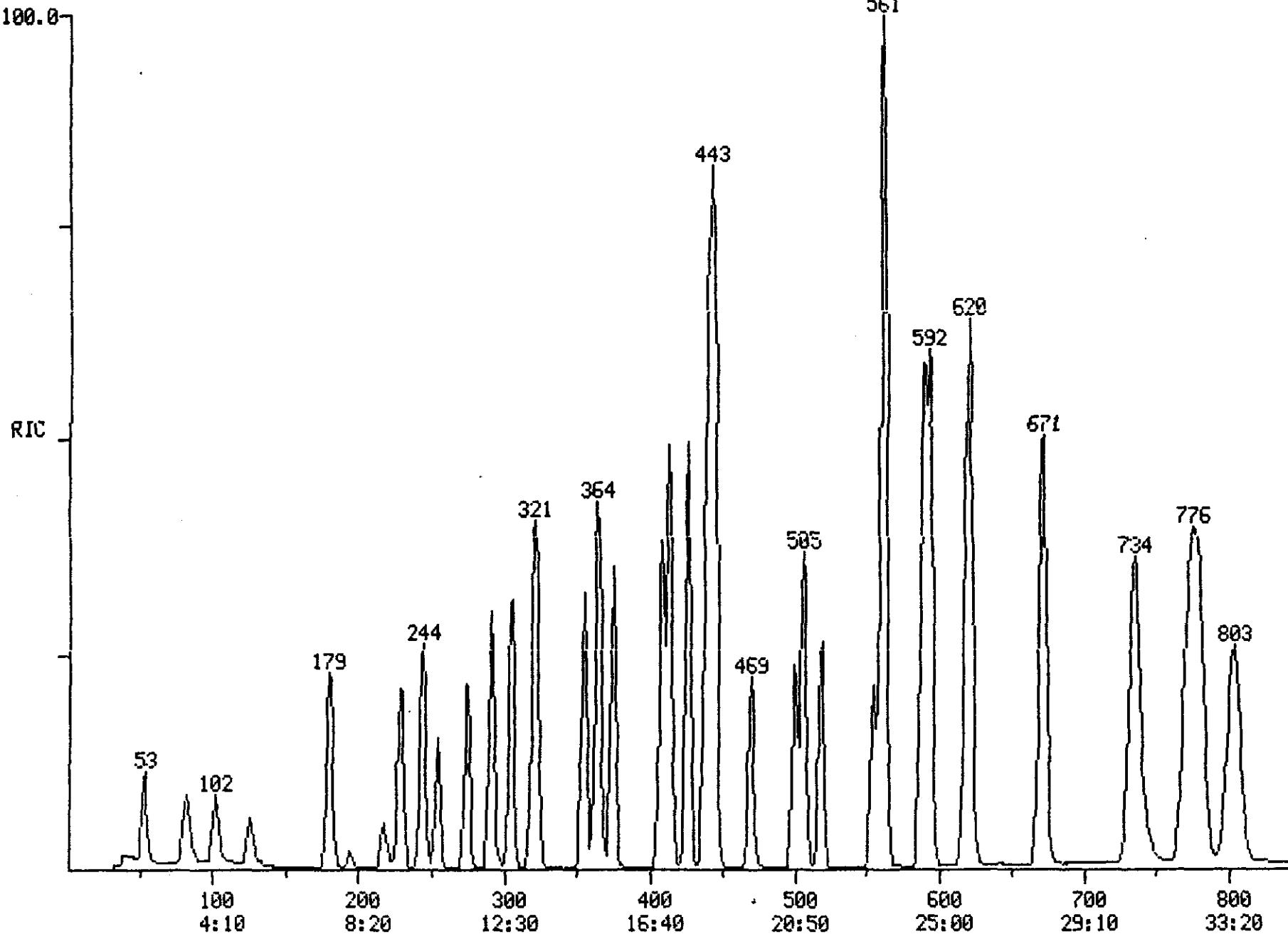
No	Ret(L)	RRT(L)	Ratio	Ret(L)	RRT(L)	Ratio	Amnt	Amnt(L)	R.Fac	R.Fac(L)	Ratio
13	13:25	1.00	1.263	1.00	1.00	1.255	1.00	50.00	50.00	1.784	1.784 1.00
14	13:20	1.00	1.255	1.00	1.00	1.255	1.00	50.00	50.00	1.790	1.790 1.00
15	9:35	1.00	0.902	1.00	1.00	0.902	1.00	50.00	50.00	1.982	1.982 1.00
16	20:47	1.00	1.000	1.00	1.00	1.000	1.00	50.00	50.00	1.000	1.000 1.00
17	13:22	1.00	1.259	1.00	1.00	1.259	1.00	50.00	50.00	0.217	0.217 1.00
18	15:15	1.00	0.733	1.00	1.00	0.733	1.00	50.00	50.00	0.558	0.558 1.00
19	15:10	1.00	0.729	1.00	1.00	0.729	1.00	50.00	50.00	0.378	0.378 1.00
20	15:37	1.00	0.752	1.00	1.00	0.752	1.00	50.00	50.00	0.511	0.511 1.00
21	17:00	1.00	0.818	1.00	1.00	0.818	1.00	50.00	50.00	0.333	0.333 1.00
22	17:12	1.00	0.828	1.00	1.00	0.828	1.00	50.00	50.00	0.600	0.600 1.00
23	17:45	1.00	0.854	1.00	1.00	0.854	1.00	50.00	50.00	0.412	0.412 1.00
24	18:17	1.00	0.880	1.00	1.00	0.880	1.00	50.00	50.00	0.827	0.827 1.00
25	18:22	1.00	0.884	1.00	1.00	0.884	1.00	50.00	50.00	0.464	0.464 1.00
26	18:30	1.00	0.890	1.00	1.00	0.890	1.00	50.00	50.00	0.315	0.315 1.00
27	18:30	1.00	0.890	1.00	1.00	0.890	1.00	50.00	50.00	0.320	0.320 1.00
28	21:02	1.00	1.012	1.00	1.00	1.012	1.00	50.00	50.00	0.433	0.433 1.00
29	14:47	1.00	0.711	1.00	1.00	0.711	1.00	50.00	50.00	0.404	0.404 1.00
30	24:30	1.00	0.953	1.00	1.00	0.953	1.00	50.00	50.00	1.213	1.213 1.00
31	25:42	1.00	1.000	1.00	1.00	1.000	1.00	50.00	50.00	1.000	1.000 1.00
32	21:35	1.00	0.840	1.00	1.00	0.840	1.00	50.00	50.00	0.652	0.652 1.00
33	23:05	1.00	0.898	1.00	1.00	0.898	1.00	50.00	50.00	0.540	0.540 1.00
34	23:22	1.00	0.909	1.00	1.00	0.909	1.00	50.00	50.00	0.399	0.399 1.00
35	23:22	1.00	0.909	1.00	1.00	0.909	1.00	50.00	50.00	0.683	0.683 1.00
36	24:42	1.00	0.961	1.00	1.00	0.961	1.00	50.00	50.00	0.621	0.621 1.00
37	25:50	1.00	1.005	1.00	1.00	1.005	1.00	50.00	50.00	0.876	0.876 1.00
38	27:55	1.00	1.086	1.00	1.00	1.086	1.00	50.00	50.00	0.396	0.396 1.00
39	32:07	1.00	1.250	1.00	1.00	1.250	1.00	50.00	50.00	0.801	0.801 1.00
40	33:27	1.00	1.301	1.00	1.00	1.301	1.00	50.00	50.00	0.448	0.448 1.00
41	32:25	1.00	1.261	1.00	1.00	1.261	1.00	50.00	50.00	0.519	0.519 1.00
42	30:35	1.00	1.190	1.00	1.00	1.190	1.00	50.00	50.00	0.832	0.832 1.00

RIC
04/23/90 18:52:00
SAMPLE: CLP,,,VSTD100,L,S,22525,V,IC-100,,,5ML,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
RANGE: G 1, 840 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: U3330 #1
CALI: U3330 #2

SCANS 1 TO 840

529408,00085



Data: U3330.TI

04/23/90 18:52:00

Sample: CLP,,, VSTD100, L, S, 22525, V, IC-100,,, 5ML,

Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

Formula: - Instrument: U Weight: 0.012

Submitted by: VERSAR Analyst: DD

Acct. No.: -

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)

Resp. fac. from Library Entry

No	Name
1	C101 BROMOCHLOROMETHANE **IS#1**
2	C010 CHLOROMETHANE
3	C015 BROMOMETHANE
4	C020 VINYL CHLORIDE
5	C025 CHLOROETHANE
6	C030 METHYLENE CHLORIDE
7	C035 ACETONE
8	C040 CARBON DISULFIDE
9	C045 1,1-DICHLOROETHENE
10	C050 1,1-DICHLOROETHANE
11	C053 1,2-DICHLOROETHENE (TOTAL)
12	C060 CHLOROFORM
13	C065 1,2-DICHLOROETHANE
14	CS15 1,2-DICHLOROETHANE-D4 **SS#1**
15	C043 TRICHLOROFLUOROMETHANE
16	C110 1,4-DIFLUOROBENZENE **IS#2**
17	C110 2-BUTANONE
18	C125 VINYL ACETATE
19	C120 CARBON TETRACHLORIDE
20	C130 BROMODICHLOROMETHANE
21	C140 1,2-DICHLOROPROPANE
22	C145 CIS-1,3-DICHLOROPROPENE
23	C150 TRICHLOROETHENE
24	C165 BENZENE
25	C155 DIBROMOCHLOROMETHANE
26	C160 1,1,2-TRICHLOROETHANE
27	C170 TRANS-1,3-DICHLOROPROPENE
28	C180 BROMOFORM
29	C115 1,1,1-TRICHLOROETHANE
30	CS05 TOLUENE-D8 **SS#2**
31	C120 CHLOROBENZENE-D5 **IS#3**
32	C205 4-METHYL-2-PENTANONE
33	C210 2-HEXANONE
34	C220 TETRACHLOROETHENE
35	C225 1,1,2,2-TETRACHLOROETHANE
36	C230 TOLUENE
37	C235 CHLOROBENZENE
38	C240 ETHYLBENZENE
39	C245 STYRENE
40	C250 ORTHO & PARA XYLENE
41	C251 META-XYLENE
42	CS10 BROMOFLUOROBENZENE **SS#3**

000087 U 3330

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	128	254	10:35	1	1.000	A BB	53689.	50.000 UG/L*	1.23
2	50	53	2:12	1	0.209	A BB	134860.	100.000 UG/L	2.47
3	94	82	3:25	1	0.323	A BB	89419.	100.000 UG/L	2.47
4	62	102	4:15	1	0.402	A BB	122456.	100.000 UG/L	2.47
5	64	126	5:15	1	0.496	A BB	65289.	100.000 UG/L	2.47
6	84	179	7:27	1	0.705	A BB	134225.	100.000 UG/L	2.47
7	43	194	8:05	1	0.764	A BB	30817.	100.000 UG/L	2.47
8	76	217	9:02	1	0.854	A BB	120245.	100.000 UG/L	2.47
9	96	244	10:10	1	0.961	A BB	124526.	100.000 UG/L	2.47
10	63	274	11:25	1	1.079	A BB	241903.	100.000 UG/L	2.47
11	96	291	12:07	1	1.146	A BB	145663.	100.000 UG/L	2.47
12	83	304	12:40	1	1.197	A BB	269138.	100.000 UG/L	2.47
13	62	322	13:25	1	1.268	A BB	189665.	100.000 UG/L	2.47
14	65	320	13:20	1	1.260	A BB	166045.	100.000 UG/L%	2.47
15	101	228	9:30	1	0.898	A BB	231085.	100.000 UG/L	2.47
16	114	499	20:47	16	1.000	A BB	242168.	50.000 UG/L*	1.23
17	72	320	13:20	1	1.260	A BB	20384.	100.000 UG/L	2.47
18	43	365	15:12	16	0.731	A BB	263466.	100.000 UG/L	2.47
19	117	363	15:07	16	0.727	A VB	217060.	100.000 UG/L	2.47
20	83	374	15:35	16	0.749	A BB	275016.	100.000 UG/L	2.47
21	63	408	17:00	16	0.818	A BB	167123.	100.000 UG/L	2.47
22	75	413	17:12	16	0.828	A BB	329532.	100.000 UG/L	2.47
23	130	426	17:45	16	0.854	A BB	216584.	100.000 UG/L	2.47
24	78	439	18:17	16	0.880	A BB	453441.	100.000 UG/L	2.47
25	129	441	18:22	16	0.884	A BB	232727.	100.000 UG/L	2.47
26	97	444	18:30	16	0.890	A BB	154332.	100.000 UG/L	2.47
27	75	443	18:27	16	0.888	A BB	175518.	100.000 UG/L	2.47
28	173	505	21:02	16	1.012	A BB	246098.	100.000 UG/L	2.47
29	97	354	14:45	16	0.709	A BB	241878.	100.000 UG/L	2.47
30	98	588	24:30	31	0.953	A BB	490301.	100.000 UG/L%	2.47
31	117	617	25:42	31	1.000	A BB	227837.	50.000 UG/L*	1.23
32	43	518	21:35	31	0.840	A BB	250526.	100.000 UG/L	2.47
33	43	554	23:05	31	0.898	A BV	211965.	100.000 UG/L	2.47
34	164	561	23:22	31	0.909	A BB	196273.	100.000 UG/L	2.47
35	83	561	23:22	31	0.909	A BB	320853.	100.000 UG/L	2.47
36	92	593	24:42	31	0.961	A BB	322912.	100.000 UG/L	2.47
37	112	620	25:50	31	1.005	A BB	410740.	100.000 UG/L	2.47
38	106	670	27:55	31	1.086	A BB	215985.	100.000 UG/L	2.47
39	104	772	32:10	31	1.251	A BB	362602.	100.000 UG/L	2.47
40	106	803	33:27	31	1.301	A BB	206056.	100.000 UG/L	2.47
41	106	778	32:25	31	1.261	A BB	238341.	100.000 UG/L	2.47
42	95	734	30:35	31	1.190	A BB	343986.	100.000 UG/L%	2.47

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
1	10:35	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
2	2:12	1.00	0.209	1.00	100.00	100.00	1.256	1.256	1.00
3	3:25	1.00	0.323	1.00	100.00	100.00	0.833	0.833	1.00
4	4:15	1.00	0.402	1.00	100.00	100.00	1.140	1.140	1.00
5	5:15	1.00	0.496	1.00	100.00	100.00	0.608	0.608	1.00
6	7:27	1.00	0.705	1.00	100.00	100.00	1.250	1.250	1.00
7	8:05	1.00	0.764	1.00	100.00	100.00	0.287	0.287	1.00
8	9:02	1.00	0.854	1.00	100.00	100.00	1.120	1.120	1.00
9	10:10	1.00	0.961	1.00	100.00	100.00	1.160	1.160	1.00
10	11:25	1.00	1.079	1.00	100.00	100.00	2.253	2.253	1.00
11	12:07	1.00	1.146	1.00	100.00	100.00	1.357	1.357	1.00
12	12:40	1.00	1.197	1.00	100.00	100.00	2.506	2.506	1.00

100087

U 333 U
10.3

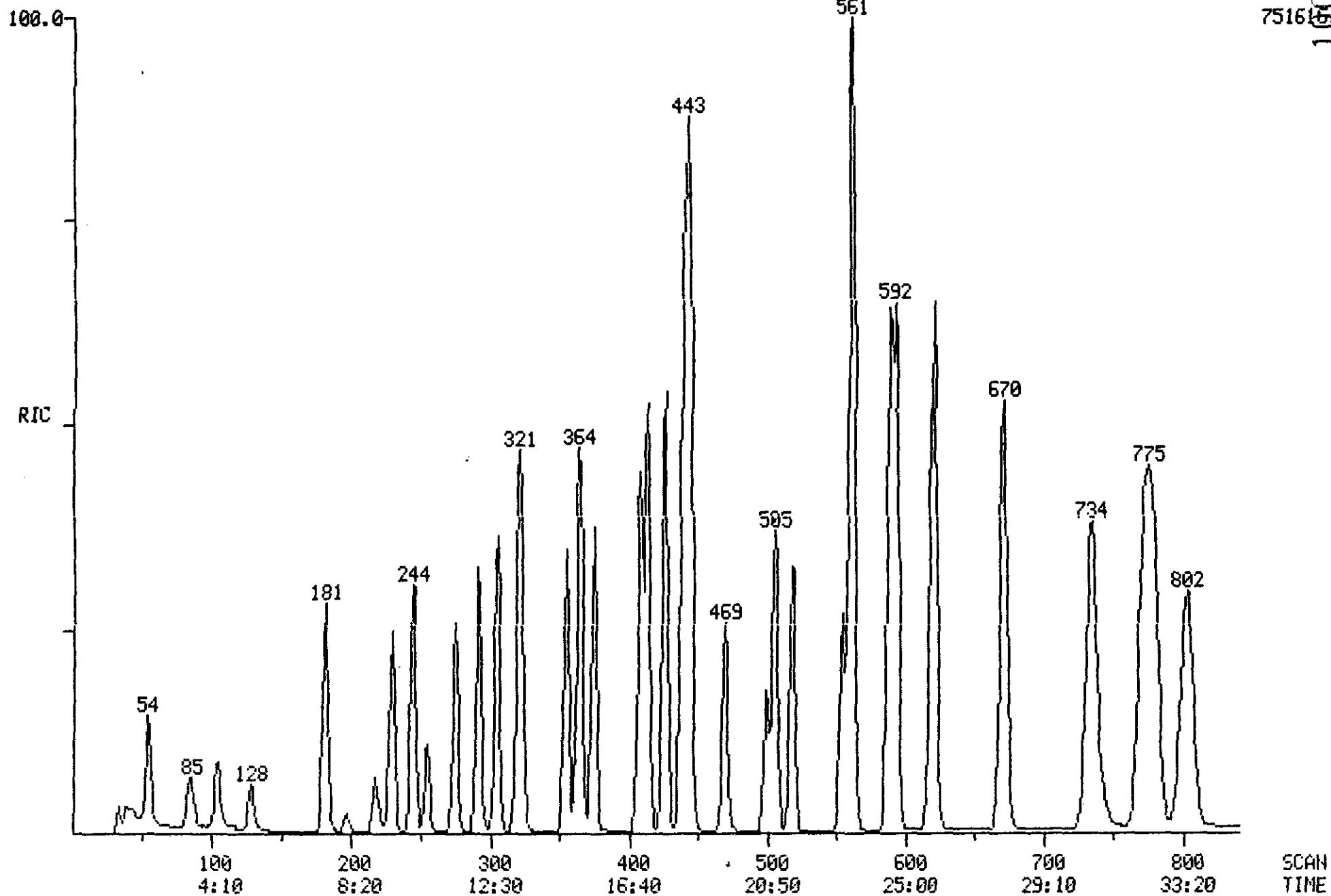
No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
13	13:25	1.00	1.268	1.00	100.00	100.00	1.766	1.766	1.00
14	13:20	1.00	1.260	1.00	100.00	100.00	1.546	1.546	1.00
15	9:30	1.00	0.898	1.00	100.00	100.00	2.152	2.152	1.00
16	20:47	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
17	13:20	1.00	1.260	1.00	100.00	100.00	0.190	0.190	1.00
18	15:12	1.00	0.731	1.00	100.00	100.00	0.544	0.544	1.00
19	15:07	1.00	0.727	1.00	100.00	100.00	0.448	0.448	1.00
20	15:35	1.00	0.749	1.00	100.00	100.00	0.568	0.568	1.00
21	17:00	1.00	0.818	1.00	100.00	100.00	0.345	0.345	1.00
22	17:12	1.00	0.828	1.00	100.00	100.00	0.680	0.680	1.00
23	17:45	1.00	0.854	1.00	100.00	100.00	0.447	0.447	1.00
24	18:17	1.00	0.880	1.00	100.00	100.00	0.936	0.936	1.00
25	18:22	1.00	0.884	1.00	100.00	100.00	0.481	0.481	1.00
26	18:30	1.00	0.890	1.00	100.00	100.00	0.319	0.319	1.00
27	18:27	1.00	0.888	1.00	100.00	100.00	0.362	0.362	1.00
28	21:02	1.00	1.012	1.00	100.00	100.00	0.508	0.508	1.00
29	14:45	1.00	0.709	1.00	100.00	100.00	0.499	0.499	1.00
30	24:30	1.00	0.953	1.00	100.00	100.00	1.076	1.076	1.00
31	25:42	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
32	21:35	1.00	0.840	1.00	100.00	100.00	0.550	0.550	1.00
33	23:05	1.00	0.898	1.00	100.00	100.00	0.465	0.465	1.00
34	23:22	1.00	0.909	1.00	100.00	100.00	0.431	0.431	1.00
35	23:22	1.00	0.909	1.00	100.00	100.00	0.704	0.704	1.00
36	24:42	1.00	0.961	1.00	100.00	100.00	0.709	0.709	1.00
37	25:50	1.00	1.005	1.00	100.00	100.00	0.901	0.901	1.00
38	27:55	1.00	1.086	1.00	100.00	100.00	0.474	0.474	1.00
39	32:10	1.00	1.251	1.00	100.00	100.00	0.796	0.796	1.00
40	33:27	1.00	1.301	1.00	100.00	100.00	0.452	0.452	1.00
41	32:25	1.00	1.261	1.00	100.00	100.00	0.523	0.523	1.00
42	30:35	1.00	1.190	1.00	100.00	100.00	0.755	0.755	1.00

RIC
04/23/90 18:09:00
SAMPLE: CLP,,,VSTD150,L,S,22524,V,IC-150,,,5ML,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ SC/MIN
RANGE: G 1, 840 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 561 3

DATA: U3329 #1
CALI: U3329 #2

SCANS 1 TO 840

75161
1600089



Quantitation Report File: U3329

SEARCHED
INDEXED
FILED

Data: U3329.TI

04/23/90 18:09:00

Sample: CLP,,, VSTD150, L, S, 22524, V, IC-150,,, 5ML,

Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

Formula: - Instrument: U Weight: 0.012

Submitted by: VERSAR Analyst: DD Acct. No.: -

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)

Resp. fac. from Library Entry

No	Name
1	C101 BROMOCHLOROMETHANE **IS#1**
2	C010 CHLOROMETHANE
3	C015 BROMOMETHANE
4	C020 VINYL CHLORIDE
5	C025 CHLOROETHANE
6	C030 METHYLENE CHLORIDE
7	C035 ACETONE
8	C040 CARBON DISULFIDE
9	C045 1,1-DICHLOROETHENE
10	C050 1,1-DICHLOROETHANE
11	C053 1,2-DICHLOROETHENE (TOTAL)
12	C060 CHLOROFORM
13	C065 1,2-DICHLOROETHANE
14	CS15 1,2-DICHLOROETHANE-D4 **SS#1**
15	C043 TRICHLOROFUOROMETHANE
16	CI10 1,4-DIFLUOROBENZENE **IS#2**
17	C110 2-BUTANONE
18	C125 VINYL ACETATE
19	C120 CARBON TETRACHLORIDE
20	C130 BROMODICHLOROMETHANE
21	C140 1,2-DICHLOROPROPANE
22	C145 CIS-1,3-DICHLOROPROPENE
23	C150 TRICHLOROETHENE
24	C165 BENZENE
25	C155 DIBROMOCHLOROMETHANE
26	C160 1,1,2-TRICHLOROETHANE
27	C170 TRANS-1,3-DICHLOROPROPENE
28	C180 BROMOFORM
29	C115 1,1,1-TRICHLOROETHANE
30	CS05 TOLUENE-D8 **SS#2**
31	CI20 CHLOROBENZENE-D5 **IS#3**
32	C205 4-METHYL-2-PENTANONE
33	C210 2-HEXANONE
34	C220 TETRACHLOROETHENE
35	C225 1,1,2,2-TETRACHLOROETHANE
36	C230 TOLUENE
37	C235 CHLOROBENZENE
38	C240 ETHYLBENZENE
39	C245 STYRENE
40	C250 ORTHO & PARA XYLENE
41	C251 META-XYLENE
42	CS10 BROMOFLUOROBENZENE **SS#3**

U3329

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	128	254	10:35	1	1.000	A BB	53747.	50.000 UG/L*	0.83
2	50	54	2:15	1	0.213	A BB	237817.	150.000 UG/L	2.50
3	94	85	3:32	1	0.335	A BB	99341.	150.000 UG/L	2.50
4	62	104	4:20	1	0.409	A BB	162659.	150.000 UG/L	2.50
5	64	128	5:20	1	0.504	A BB	93106.	150.000 UG/L	2.50
6	84	181	7:32	1	0.713	A BB	212375.	150.000 UG/L	2.50
7	43	197	8:12	1	0.776	A BB	54860.	150.000 UG/L	2.50
8	76	218	9:05	1	0.858	A BB	200009.	150.000 UG/L	2.50
9	96	244	10:10	1	0.961	A BB	196958.	150.000 UG/L	2.50
10	63	275	11:27	1	1.083	A BB	389207.	150.000 UG/L	2.50
11	96	291	12:07	1	1.146	A BB	216596.	150.000 UG/L	2.50
12	83	305	12:42	1	1.201	A BV	420054.	150.000 UG/L	2.50
13	62	322	13:25	1	1.268	A BB	298481.	150.000 UG/L	2.50
14	65	320	13:20	1	1.260	A BB	263071.	150.000 UG/L%	2.50
15	101	229	9:32	1	0.902	A BV	361755.	150.000 UG/L	2.50
16	114	499	20:47	16	1.000	A BB	244897.	50.000 UG/L*	0.83
17	72	321	13:22	1	1.264	A BB	34754.	150.000 UG/L	2.50
18	43	365	15:12	16	0.731	A BB	450762.	150.000 UG/L	2.50
19	117	364	15:10	16	0.729	A VB	324328.	150.000 UG/L	2.50
20	83	374	15:35	16	0.749	A BB	403995.	150.000 UG/L	2.50
21	63	408	17:00	16	0.818	A BB	269397.	150.000 UG/L	2.50
22	75	413	17:12	16	0.828	A BB	483101.	150.000 UG/L	2.50
23	130	426	17:45	16	0.854	A BV	330272.	150.000 UG/L	2.50
24	78	438	18:15	16	0.878	A BB	663409.	150.000 UG/L	2.50
25	129	441	18:22	16	0.884	A BB	334087.	150.000 UG/L	2.50
26	97	443	18:27	16	0.888	A BB	236773.	150.000 UG/L	2.50
27	75	443	18:27	16	0.888	A BB	254312.	150.000 UG/L	2.50
28	173	505	21:02	16	1.012	A BB	349604.	150.000 UG/L	2.50
29	97	355	14:47	16	0.711	A BB	354517.	150.000 UG/L	2.50
30	98	588	24:30	31	0.953	A BB	733141.	150.000 UG/L%	2.50
31	117	617	25:42	31	1.000	A BB	225904.	50.000 UG/L*	0.83
32	43	518	21:35	31	0.840	A BB	450561.	150.000 UG/L	2.50
33	43	554	23:05	31	0.898	A BV	382905.	150.000 UG/L	2.50
34	164	561	23:22	31	0.909	A BB	282622.	150.000 UG/L	2.50
35	83	560	23:20	31	0.908	A BB	473706.	150.000 UG/L	2.50
36	92	592	24:40	31	0.959	A BB	476505.	150.000 UG/L	2.50
37	112	620	25:50	31	1.005	A BB	624858.	150.000 UG/L	2.50
38	106	670	27:55	31	1.086	A BB	318425.	150.000 UG/L	2.50
39	104	771	32:07	31	1.250	A BB	585159.	150.000 UG/L	2.50
40	106	802	33:25	31	1.300	A BB	334825.	150.000 UG/L	2.50
41	106	778	32:25	31	1.261	A BB	383559.	150.000 UG/L	2.50
42	95	734	30:35	31	1.190	A BB	520031.	150.000 UG/L%	2.50

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
1	10:35	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
2	2:15	1.00	0.213	1.00	150.00	150.00	1.475	1.475	1.00
3	3:32	1.00	0.335	1.00	150.00	150.00	0.616	0.616	1.00
4	4:20	1.00	0.409	1.00	150.00	150.00	1.009	1.009	1.00
5	5:20	1.00	0.504	1.00	150.00	150.00	0.577	0.577	1.00
6	7:32	1.00	0.713	1.00	150.00	150.00	1.317	1.317	1.00
7	8:12	1.00	0.776	1.00	150.00	150.00	0.340	0.340	1.00
8	9:05	1.00	0.858	1.00	150.00	150.00	1.240	1.240	1.00
9	10:10	1.00	0.961	1.00	150.00	150.00	1.222	1.222	1.00
10	11:27	1.00	1.083	1.00	150.00	150.00	2.414	2.414	1.00
11	12:07	1.00	1.146	1.00	150.00	150.00	1.343	1.343	1.00
12	12:42	1.00	1.201	1.00	150.00	150.00	2.605	2.605	1.00

100091

U3291

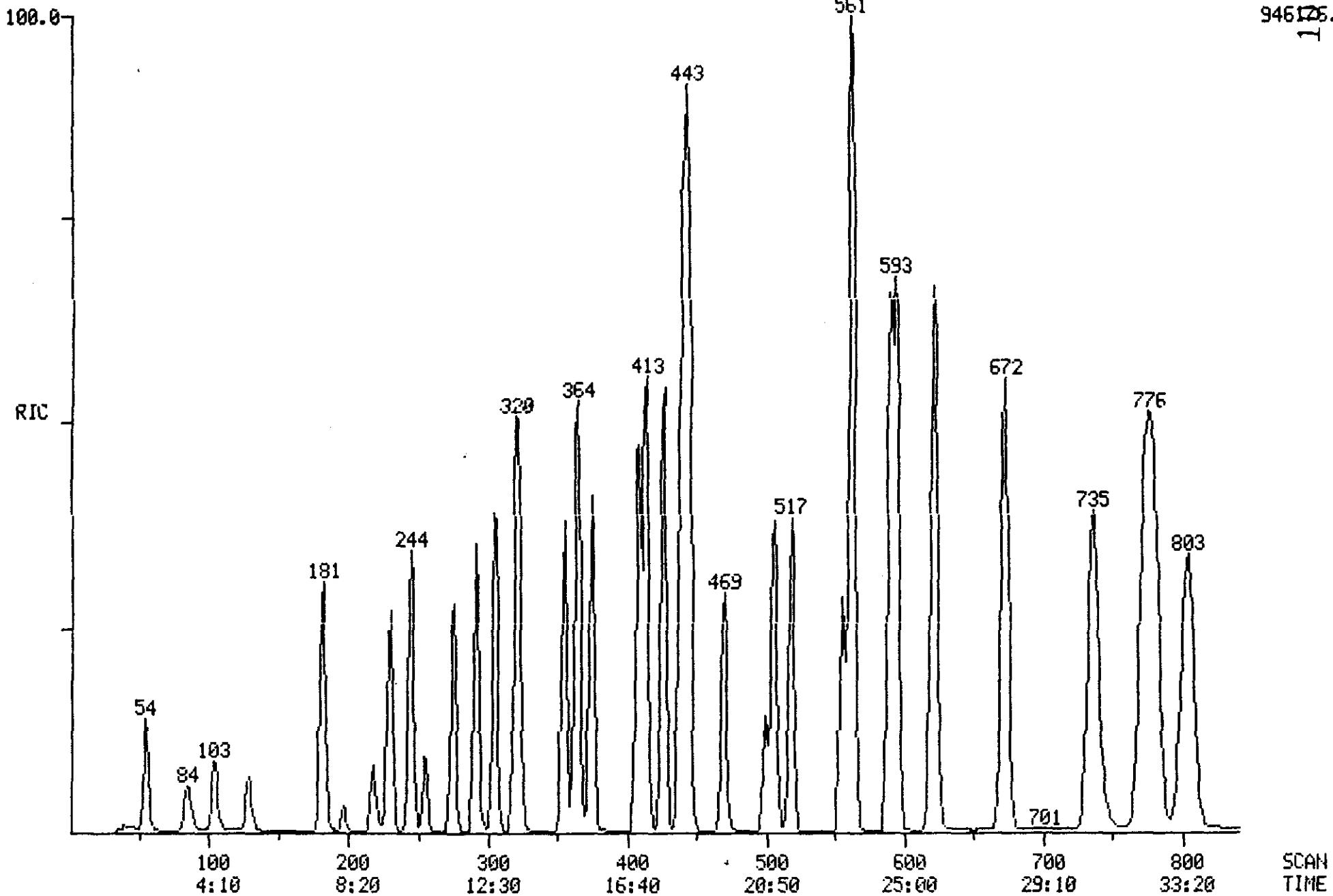
No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
13	13: 25	1.00	1. 268	1.00	150.00	150.00	1. 851	1. 851	1.00
14	13: 20	1.00	1. 260	1.00	150.00	150.00	1. 632	1. 632	1.00
15	9: 32	1.00	0. 902	1.00	150.00	150.00	2. 244	2. 244	1.00
16	20: 47	1.00	1. 000	1.00	50.00	50.00	1. 000	1. 000	1.00
17	13: 22	1.00	1. 264	1.00	150.00	150.00	0. 216	0. 216	1.00
18	15: 12	1.00	0. 731	1.00	150.00	150.00	0. 614	0. 614	1.00
19	15: 10	1.00	0. 729	1.00	150.00	150.00	0. 441	0. 441	1.00
20	15: 35	1.00	0. 749	1.00	150.00	150.00	0. 550	0. 550	1.00
21	17: 00	1.00	0. 818	1.00	150.00	150.00	0. 367	0. 367	1.00
22	17: 12	1.00	0. 828	1.00	150.00	150.00	0. 658	0. 658	1.00
23	17: 45	1.00	0. 854	1.00	150.00	150.00	0. 450	0. 450	1.00
24	18: 15	1.00	0. 878	1.00	150.00	150.00	0. 903	0. 903	1.00
25	18: 22	1.00	0. 884	1.00	150.00	150.00	0. 455	0. 455	1.00
26	18: 27	1.00	0. 888	1.00	150.00	150.00	0. 322	0. 322	1.00
27	18: 27	1.00	0. 888	1.00	150.00	150.00	0. 346	0. 346	1.00
28	21: 02	1.00	1. 012	1.00	150.00	150.00	0. 476	0. 476	1.00
29	14: 47	1.00	0. 711	1.00	150.00	150.00	0. 483	0. 483	1.00
30	24: 30	1.00	0. 953	1.00	150.00	150.00	1. 082	1. 082	1.00
31	25: 42	1.00	1. 000	1.00	50.00	50.00	1. 000	1. 000	1.00
32	21: 35	1.00	0. 840	1.00	150.00	150.00	0. 665	0. 665	1.00
33	23: 05	1.00	0. 898	1.00	150.00	150.00	0. 565	0. 565	1.00
34	23: 22	1.00	0. 909	1.00	150.00	150.00	0. 417	0. 417	1.00
35	23: 20	1.00	0. 908	1.00	150.00	150.00	0. 699	0. 699	1.00
36	24: 40	1.00	0. 959	1.00	150.00	150.00	0. 703	0. 703	1.00
37	25: 50	1.00	1. 005	1.00	150.00	150.00	0. 922	0. 922	1.00
38	27: 55	1.00	1. 086	1.00	150.00	150.00	0. 470	0. 470	1.00
39	32: 07	1.00	1. 250	1.00	150.00	150.00	0. 863	0. 863	1.00
40	33: 25	1.00	1. 300	1.00	150.00	150.00	0. 497	0. 497	1.00
41	32: 25	1.00	1. 261	1.00	150.00	150.00	0. 566	0. 566	1.00
42	30: 35	1.00	1. 190	1.00	150.00	150.00	0. 767	0. 767	1.00

RIC
04/23/90 17:21:00
SAMPLE: CLP,,,VSTD200,L,S,22523,U,IC-200,,,5ML,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
RANGE: G 1, 840 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: U3328 #1
CALI: U3328 #2

SCANS 1 TO 840

1000093
946125.



Data: U3328.TI

04/23/90 17:21:00

Sample: CLP,,,VSTD200,L,S,22523,V,IC-200,,,5ML,

Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

Formula: - Instrument: U Weight: 0.012

Submitted by: VERSAR Analyst: DD Acct. No.: -

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)

Resp. fac. from Library Entry

No	Name
1	C101 BROMOCHLOROMETHANE **IS#1**
2	C010 CHLOROMETHANE
3	C015 BROMOMETHANE
4	C020 VINYL CHLORIDE
5	C025 CHLOROETHANE
6	C030 METHYLENE CHLORIDE
7	C035 ACETONE
8	C040 CARBON DISULFIDE
9	C045 1,1-DICHLOROETHENE
10	C050 1,1-DICHLOROETHANE
11	C053 1,2-DICHLOROETHENE (TOTAL)
12	C060 CHLOROFORM
13	C065 1,2-DICHLOROETHANE
14	CS15 1,2-DICHLOROETHANE-D4 **SS#1**
15	C043 TRICHLOROFLUOROMETHANE
16	CI10 1,4-DIFLUOROBENZENE **IS#2**
17	C110 2-BUTANONE
18	C125 VINYL ACETATE
19	C120 CARBON TETRACHLORIDE
20	C130 BROMODICHLOROMETHANE
21	C140 1,2-DICHLOROPROPANE
22	C145 CIS-1,3-DICHLOROPROPENE
23	C150 TRICHLOROETHENE
24	C165 BENZENE
25	C155 DIBROMOCHLOROMETHANE
26	C160 1,1,2-TRICHLOROETHANE
27	C170 TRANS-1,3-DICHLOROPROPENE
28	C180 BROMOFORM
29	C115 1,1,1-TRICHLOROETHANE
30	CS05 TOLUENE-D8 **SS#2**
31	CI20 CHLOROBENZENE-D5 **IS#3**
32	C205 4-METHYL-2-PENTANONE
33	C210 2-HEXANONE
34	C220 TETRACHLOROETHENE
35	C225 1,1,2,2-TETRACHLOROETHANE
36	C230 TOLUENE
37	C235 CHLOROBENZENE
38	C240 ETHYLBENZENE
39	C245 STYRENE
40	C250 ORTHO & PARA XYLENE
41	C251 META-XYLENE
42	CS10 BROMOFLUOROBENZENE **SS#3**

W3328

(Ref)

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	128	254	10:35	1	1.000	A BB	54651.	50.000 UG/L*	0.63
2	50	54	2:15	1	0.213	A BB	297663.	200.000 UG/L	2.52
3	94	84	3:30	1	0.331	A BB	109635.	200.000 UG/L	2.52
4	62	103	4:17	1	0.406	A BB	212773.	200.000 UG/L	2.52
5	64	128	5:20	1	0.504	A BB	132480.	200.000 UG/L	2.52
6	84	181	7:32	1	0.713	A BB	284031.	200.000 UG/L	2.52
7	43	196	8:10	1	0.772	A BB	76944.	200.000 UG/L	2.52
8	76	217	9:02	1	0.854	A BB	285027.	200.000 UG/L	2.52
9	96	244	10:10	1	0.961	A BB	264863.	200.000 UG/L	2.52
10	63	275	11:27	1	1.083	A BB	534173.	200.000 UG/L	2.52
11	96	291	12:07	1	1.146	A BB	290009.	200.000 UG/L	2.52
12	83	304	12:40	1	1.197	A BV	564119.	200.000 UG/L	2.52
13	62	322	13:25	1	1.268	A BB	415017.	200.000 UG/L	2.52
14	65	319	13:17	1	1.256	A BB	360329.	200.000 UG/L%	2.52
15	101	229	9:32	1	0.902	A BV	487557.	200.000 UG/L	2.52
16	114	499	20:47	16	1.000	A BB	246636.	50.000 UG/L*	0.63
17	72	320	13:20	1	1.260	A BB	47863.	200.000 UG/L	2.52
18	43	365	15:12	16	0.731	A BB	676553.	200.000 UG/L	2.52
19	117	363	15:07	16	0.727	A VB	423736.	200.000 UG/L	2.52
20	83	374	15:35	16	0.749	A BB	542994.	200.000 UG/L	2.52
21	63	407	16:57	16	0.816	A BB	361877.	200.000 UG/L	2.52
22	75	413	17:12	16	0.828	A BB	644101.	200.000 UG/L	2.52
23	130	425	17:42	16	0.852	A BV	424229.	200.000 UG/L	2.52
24	78	438	18:15	16	0.878	A BB	874082.	200.000 UG/L	2.52
25	129	440	18:20	16	0.882	A BB	418000.	200.000 UG/L	2.52
26	97	443	18:27	16	0.888	A BB	309621.	200.000 UG/L	2.52
27	75	443	18:27	16	0.888	A BB	341668.	200.000 UG/L	2.52
28	173	505	21:02	16	1.012	A BB	447819.	200.000 UG/L	2.52
29	97	354	14:45	16	0.709	A BB	473049.	200.000 UG/L	2.52
30	98	589	24:32	31	0.953	A BB	934041.	200.000 UG/L%	2.52
31	117	618	25:45	31	1.000	A BB	229121.	50.000 UG/L*	0.63
32	43	517	21:32	31	0.837	A BB	662164.	200.000 UG/L	2.52
33	43	555	23:07	31	0.898	A BB	519484.	200.000 UG/L	2.52
34	164	561	23:22	31	0.908	A BB	353490.	200.000 UG/L	2.52
35	83	561	23:22	31	0.908	A BB	618932.	200.000 UG/L	2.52
36	92	593	24:42	31	0.960	A BB	620898.	200.000 UG/L	2.52
37	112	621	25:52	31	1.005	A BB	817646.	200.000 UG/L	2.52
38	106	671	27:57	31	1.086	A BB	414518.	200.000 UG/L	2.52
39	104	772	32:10	31	1.249	A BB	828038.	200.000 UG/L	2.52
40	106	803	33:27	31	1.299	A BB	480534.	200.000 UG/L	2.52
41	106	779	32:27	31	1.261	A BB	548884.	200.000 UG/L	2.52
42	95	735	30:37	31	1.189	A BB	670094.	200.000 UG/L%	2.52

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R.Fac	R.Fac(L)	Ratio
1	10:35	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
2	2:15	1.00	0.213	1.00	200.00	200.00	1.362	1.362	1.00
3	3:30	1.00	0.331	1.00	200.00	200.00	0.502	0.502	1.00
4	4:17	1.00	0.406	1.00	200.00	200.00	0.973	0.973	1.00
5	5:20	1.00	0.504	1.00	200.00	200.00	0.606	0.606	1.00
6	7:32	1.00	0.713	1.00	200.00	200.00	1.299	1.299	1.00
7	8:10	1.00	0.772	1.00	200.00	200.00	0.352	0.352	1.00
8	9:02	1.00	0.854	1.00	200.00	200.00	1.304	1.304	1.00
9	10:10	1.00	0.961	1.00	200.00	200.00	1.212	1.212	1.00
10	11:27	1.00	1.083	1.00	200.00	200.00	2.444	2.444	1.00
11	12:07	1.00	1.146	1.00	200.00	200.00	1.327	1.327	1.00
12	12:40	1.00	1.197	1.00	200.00	200.00	2.581	2.581	1.00

100095

ORIGINAL
(Rev)

U3328

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
13	13:25	1.00	1.268	1.00	200.00	200.00	1.898	1.898	1.00
14	13:17	1.00	1.256	1.00	200.00	200.00	1.648	1.648	1.00
15	9:32	1.00	0.902	1.00	200.00	200.00	2.230	2.230	1.00
16	20:47	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
17	13:20	1.00	1.260	1.00	200.00	200.00	0.219	0.219	1.00
18	15:12	1.00	0.731	1.00	200.00	200.00	0.686	0.686	1.00
19	15:07	1.00	0.727	1.00	200.00	200.00	0.430	0.430	1.00
20	15:35	1.00	0.749	1.00	200.00	200.00	0.550	0.550	1.00
21	16:57	1.00	0.816	1.00	200.00	200.00	0.367	0.367	1.00
22	17:12	1.00	0.828	1.00	200.00	200.00	0.653	0.653	1.00
23	17:42	1.00	0.852	1.00	200.00	200.00	0.430	0.430	1.00
24	18:15	1.00	0.878	1.00	200.00	200.00	0.886	0.886	1.00
25	18:20	1.00	0.882	1.00	200.00	200.00	0.424	0.424	1.00
26	18:27	1.00	0.888	1.00	200.00	200.00	0.314	0.314	1.00
27	18:27	1.00	0.888	1.00	200.00	200.00	0.346	0.346	1.00
28	21:02	1.00	1.012	1.00	200.00	200.00	0.454	0.454	1.00
29	14:45	1.00	0.709	1.00	200.00	200.00	0.480	0.480	1.00
30	24:32	1.00	0.953	1.00	200.00	200.00	1.019	1.019	1.00
31	25:45	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
32	21:32	1.00	0.837	1.00	200.00	200.00	0.723	0.723	1.00
33	23:07	1.00	0.898	1.00	200.00	200.00	0.567	0.567	1.00
34	23:22	1.00	0.908	1.00	200.00	200.00	0.386	0.386	1.00
35	23:22	1.00	0.908	1.00	200.00	200.00	0.675	0.675	1.00
36	24:42	1.00	0.960	1.00	200.00	200.00	0.677	0.677	1.00
37	25:52	1.00	1.005	1.00	200.00	200.00	0.452	0.452	1.00
38	27:57	1.00	1.086	1.00	200.00	200.00	0.903	0.903	1.00
39	32:10	1.00	1.249	1.00	200.00	200.00	0.524	0.524	1.00
40	33:27	1.00	1.299	1.00	200.00	200.00	0.599	0.599	1.00
41	32:27	1.00	1.261	1.00	200.00	200.00	0.731	0.731	1.00
42	30:37	1.00	1.189	1.00	200.00	200.00			

100096

RIC

04/26/90 8:30:00

SAMPLE: CLP,,,USTD50,L,S,22546,U,CC-050,,5 ML,,

COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

RANGE: G 1, 840 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: U3386 #1

CALI: U3386 #2

SCANS 1 TO 840

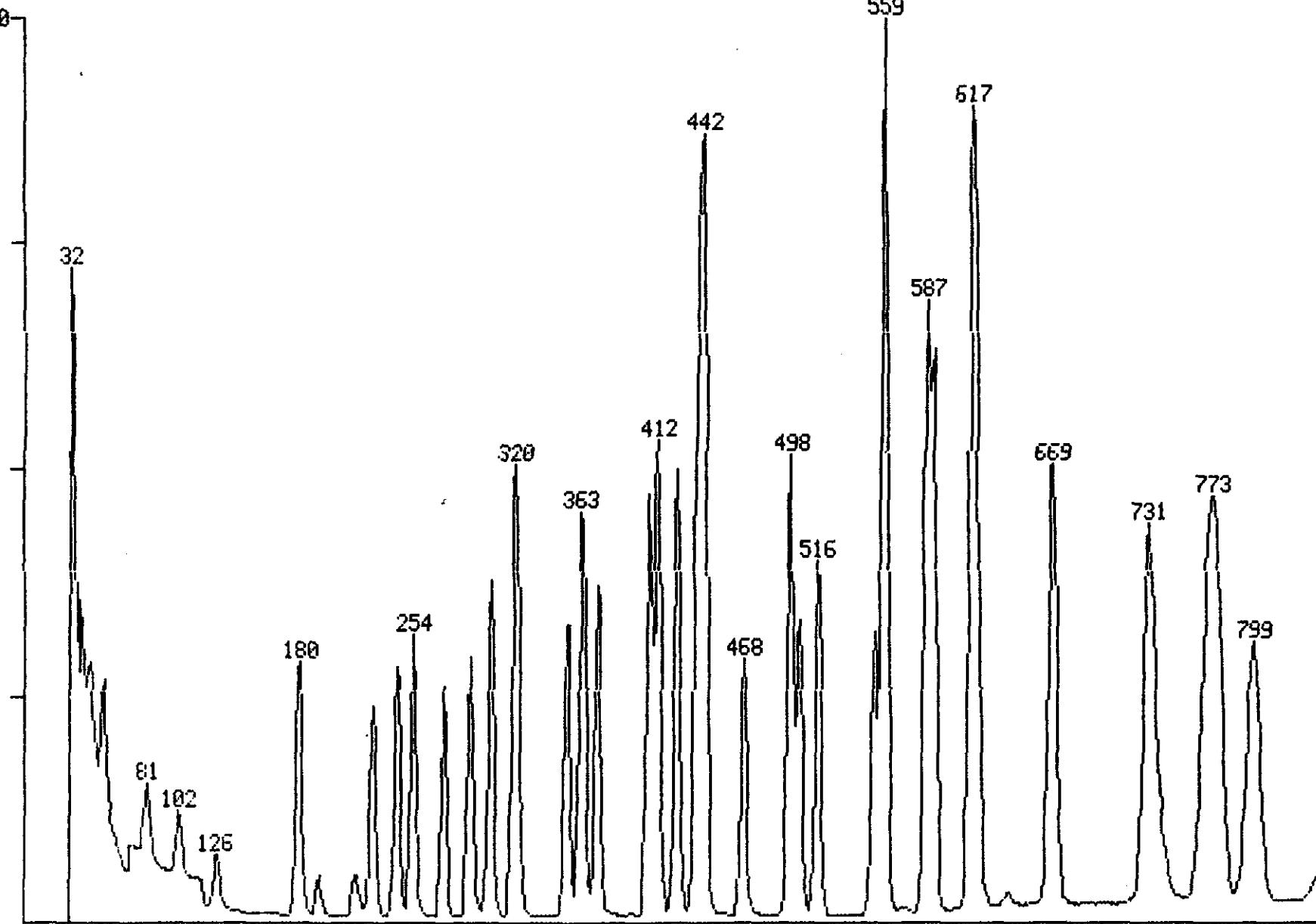
3100003?

314296.

100.0

RIC

32

100
4:10200
8:20300
12:30400
16:40500
20:50600
25:00700
29:10800
33:20SCAN
TIME

Quantitation Report File: U3386

6/10/91
(H&P)

Data: U3386.TI

04/26/90 8:30:00

Sample: CLP,,, VSTD50, L, S, 22546, V, CC-050,, 5 ML.,

Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

Formula: -

Instrument: U

Weight: 0.011

Submitted by: VERSAR

Analyst: MAT

Acct. No.: -

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)

Resp. fac. from Library Entry

No	Name
1	C101 BROMOCHLOROMETHANE **IS#1**
2	C010 CHLOROMETHANE
3	C015 BROMOMETHANE
4	C020 VINYL CHLORIDE
5	C025 CHLOROETHANE
6	C030 METHYLENE CHLORIDE
7	C035 ACETONE
8	C040 CARBON DISULFIDE
9	C045 1,1-DICHLOROETHENE
10	C050 1,1-DICHLOROETHANE
11	C053 1,2-DICHLOROETHENE (TOTAL)
12	C060 CHLOROFORM
13	C065 1,2-DICHLOROETHANE
14	CS15 1,2-DICHLOROETHANE-D4 **SS#1**
15	C043 TRICHLOROFLUOROMETHANE
16	CI10 1,4-DIFLUOROBENZENE **IS#2**
17	C110 2-BUTANONE
18	C125 VINYL ACETATE
19	C120 CARBON TETRACHLORIDE
20	C130 BROMODICHLOROMETHANE
21	C140 1,2-DICHLOROPROPANE
22	C145 CIS-1,3-DICHLOROPROPENE
23	C150 TRICHLOROETHENE
24	C165 BENZENE
25	C155 DIBROMOCHLOROMETHANE
26	C160 1,1,2-TRICHLOROETHANE
27	C170 TRANS-1,3-DICHLOROPROPENE
28	C180 BROMOFORM
29	C115 1,1,1-TRICHLOROETHANE
30	CS05 TOLUENE-D8 **SS#2**
31	CI20 CHLOROBENZENE-D5 **IS#3**
32	C205 4-METHYL-2-PENTANONE
33	C210 2-HEXANONE
34	C220 TETRACHLOROETHENE
35	C225 1,1,2,2-TETRACHLOROETHANE
36	C230 TOLUENE
37	C235 CHLOROBENZENE
38	C240 ETHYLBENZENE
39	C245 STYRENE
40	C250 ORTHO & PARA XYLENE
41	C251 META-XYLENE
42	CS10 BROMOFLUOROBENZENE **SS#3**

100002

43386

ORIGINAL

100099

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	128	253	10:32	1	1.000	A BB	55811.	50.000 UG/L*	2.38
2	50	53	2:12	1	0.209	A BB	75255.	50.000 UG/L	2.38
3	94	81	3:22	1	0.320	A BB	47726.	50.000 UG/L	2.38
4	62	102	4:15	1	0.403	A BB	56898.	50.000 UG/L	2.38
5	64	126	5:15	1	0.498	A BB	38043.	50.000 UG/L	2.38
6	84	179	7:27	1	0.708	A BB	76571.	50.000 UG/L	2.38
7	43	193	8:02	1	0.763	A BB	40566.	50.000 UG/L	2.38
8	76	217	9:02	1	0.858	A BB	58305.	50.000 UG/L	2.38
9	96	243	10:07	1	0.960	A BB	64277.	50.000 UG/L	2.38
10	63	274	11:25	1	1.083	A BB	158324.	50.000 UG/L	2.38
11	96	290	12:05	1	1.146	A BB	75558.	50.000 UG/L	2.38
12	83	304	12:40	1	1.202	A BB	165158.	50.000 UG/L	2.38
13	62	321	13:22	1	1.269	A BB	140234.	50.000 UG/L	2.38
14	65	319	13:17	1	1.261	A BB	127679.	50.000 UG/L%	2.38
15	101	228	9:30	1	0.901	A BB	129320.	50.000 UG/L	2.38
16	114	498	20:45	16	1.000	A BB	266706.	50.000 UG/L*	2.38
17	72	319	13:17	1	1.261	A BB	13148.	50.000 UG/L	2.38
18	43	364	15:10	16	0.731	A BB	210945.	50.000 UG/L	2.38
19	117	363	15:07	16	0.729	A VB	114071.	50.000 UG/L	2.38
20	83	373	15:32	16	0.749	A BB	154783.	50.000 UG/L	2.38
21	63	407	16:57	16	0.817	A BB	104618.	50.000 UG/L	2.38
22	75	412	17:10	16	0.827	A BB	180994.	50.000 UG/L	2.38
23	130	425	17:42	16	0.853	A BB	115430.	50.000 UG/L	2.38
24	78	437	18:12	16	0.878	A BB	237225.	50.000 UG/L	2.38
25	129	440	18:20	16	0.884	A BB	129732.	50.000 UG/L	2.38
26	97	442	18:25	16	0.888	A BB	93916.	50.000 UG/L	2.38
27	75	442	18:25	16	0.888	A BB	97645.	50.000 UG/L	2.38
28	173	504	21:00	16	1.012	A BB	118876.	50.000 UG/L	2.38
29	97	353	14:42	16	0.709	A BB	131529.	50.000 UG/L	2.38
30	98	587	24:27	31	0.953	A BB	304802.	50.000 UG/L%	2.38
31	117	616	25:40	31	1.000	A BB	253527.	50.000 UG/L*	2.38
32	43	516	21:30	31	0.838	A BB	233709.	50.000 UG/L	2.38
33	43	553	23:02	31	0.898	A BB	192526.	50.000 UG/L	2.38
34	164	559	23:17	31	0.907	A BB	105650.	50.000 UG/L	2.38
35	83	559	23:17	31	0.907	A BB	180766.	50.000 UG/L	2.38
36	92	591	24:37	31	0.959	A BB	174006.	50.000 UG/L	2.38
37	112	619	25:47	31	1.005	A BB	231882.	50.000 UG/L	2.38
38	106	668	27:50	31	1.084	A BB	111260.	50.000 UG/L	2.38
39	104	768	32:00	31	1.247	A BB	218497.	50.000 UG/L	2.38
40	106	799	33:17	31	1.297	A BB	120984.	50.000 UG/L	2.38
41	106	774	32:15	31	1.256	A BB	141730.	50.000 UG/L	2.38
42	95	731	30:27	31	1.187	A BB	220308.	50.000 UG/L%	2.38

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R.Fac	R.Fac(L)	Ratio
1	10:32	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
2	2:12	1.00	0.209	1.00	50.00	50.00	1.348	1.348	1.00
3	3:22	1.00	0.320	1.00	50.00	50.00	0.855	0.855	1.00
4	4:15	1.00	0.403	1.00	50.00	50.00	1.019	1.019	1.00
5	5:15	1.00	0.498	1.00	50.00	50.00	0.682	0.682	1.00
6	7:27	1.00	0.708	1.00	50.00	50.00	1.372	1.372	1.00
7	8:02	1.00	0.763	1.00	50.00	50.00	0.727	0.727	1.00
8	9:02	1.00	0.858	1.00	50.00	50.00	1.045	1.045	1.00
9	10:07	1.00	0.960	1.00	50.00	50.00	1.152	1.152	1.00
10	11:25	1.00	1.083	1.00	50.00	50.00	2.837	2.837	1.00
11	12:05	1.00	1.146	1.00	50.00	50.00	1.354	1.354	1.00
12	12:40	1.00	1.202	1.00	50.00	50.00	2.959	2.959	1.00

100099

U 3386

JUN 1986

(Ref)

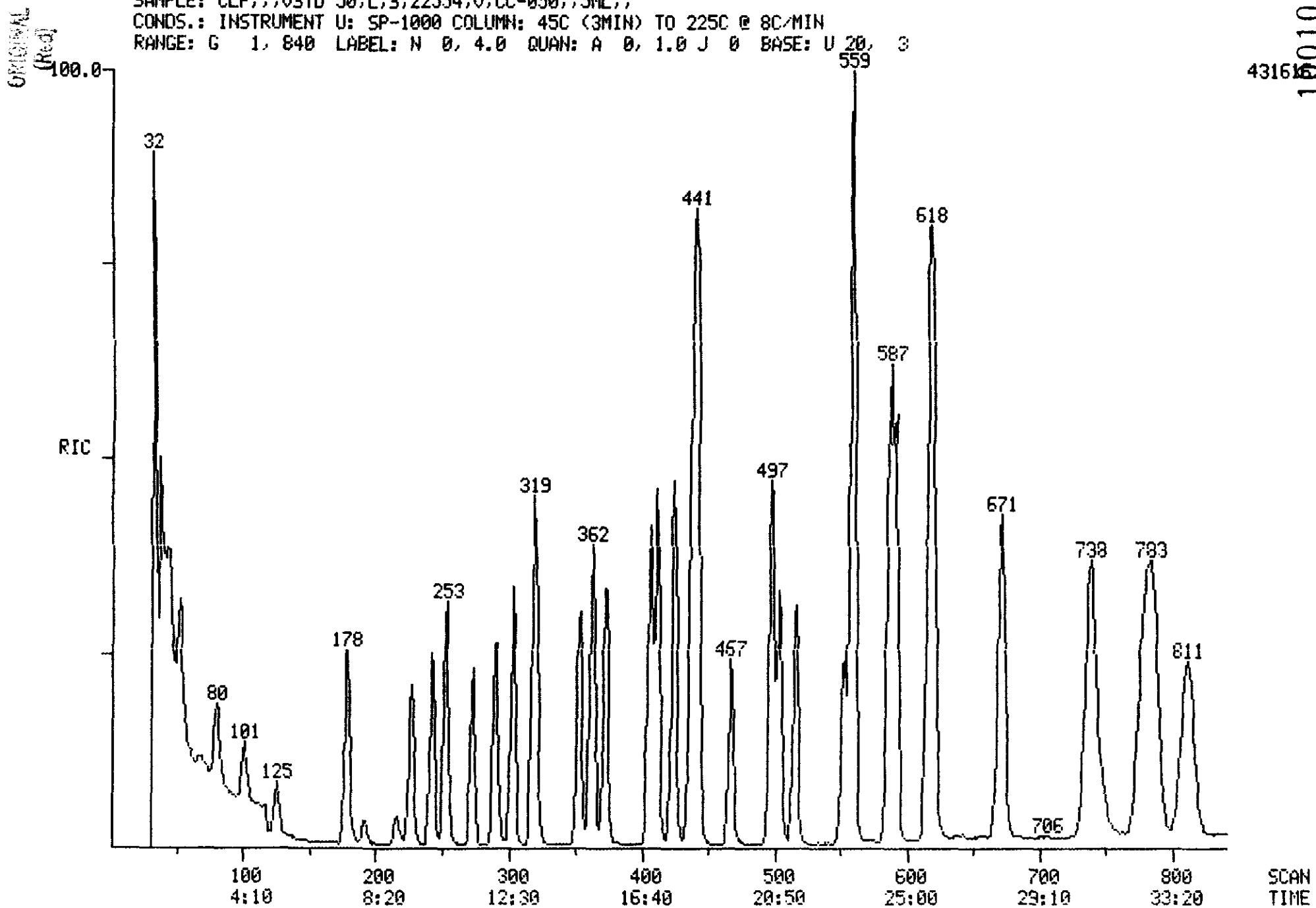
No	Ret(L)	RRT(L)	Ratio	Ret(L)	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
13	13:22	1.00	1.269	1.00	1.00	1.269	50.00	50.00	2.513	2.513	1.00
14	13:17	1.00	1.261	1.00	1.00	1.261	50.00	50.00	2.288	2.288	1.00
15	9:30	1.00	0.901	1.00	1.00	0.901	50.00	50.00	2.317	2.317	1.00
16	20:45	1.00	1.000	1.00	1.00	1.000	50.00	50.00	1.000	1.000	1.00
17	13:17	1.00	1.261	1.00	1.00	1.261	50.00	50.00	0.236	0.236	1.00
18	15:10	1.00	0.731	1.00	1.00	0.731	50.00	50.00	0.791	0.791	1.00
19	15:07	1.00	0.729	1.00	1.00	0.729	50.00	50.00	0.428	0.428	1.00
20	20:15	1.00	0.749	1.00	1.00	0.749	50.00	50.00	0.580	0.580	1.00
21	16:57	1.00	0.817	1.00	1.00	0.817	50.00	50.00	0.392	0.392	1.00
22	17:10	1.00	0.827	1.00	1.00	0.827	50.00	50.00	0.679	0.679	1.00
23	17:42	1.00	0.853	1.00	1.00	0.853	50.00	50.00	0.433	0.433	1.00
24	18:12	1.00	0.878	1.00	1.00	0.878	50.00	50.00	0.889	0.889	1.00
25	18:20	1.00	0.884	1.00	1.00	0.884	50.00	50.00	0.486	0.486	1.00
26	18:25	1.00	0.888	1.00	1.00	0.888	50.00	50.00	0.352	0.352	1.00
27	18:25	1.00	0.888	1.00	1.00	0.888	50.00	50.00	0.366	0.366	1.00
28	21:00	1.00	1.012	1.00	1.00	1.012	50.00	50.00	0.446	0.446	1.00
29	14:42	1.00	0.709	1.00	1.00	0.709	50.00	50.00	0.493	0.493	1.00
30	24:27	1.00	0.953	1.00	1.00	0.953	50.00	50.00	1.202	1.202	1.00
31	25:40	1.00	1.000	1.00	1.00	1.000	50.00	50.00	1.000	1.000	1.00
32	21:30	1.00	0.838	1.00	1.00	0.838	50.00	50.00	0.922	0.922	1.00
33	23:02	1.00	0.898	1.00	1.00	0.898	50.00	50.00	0.759	0.759	1.00
34	23:17	1.00	0.907	1.00	1.00	0.907	50.00	50.00	0.417	0.417	1.00
35	23:17	1.00	0.907	1.00	1.00	0.907	50.00	50.00	0.713	0.713	1.00
36	24:37	1.00	0.959	1.00	1.00	0.959	50.00	50.00	0.686	0.686	1.00
37	25:47	1.00	1.005	1.00	1.00	1.005	50.00	50.00	0.915	0.915	1.00
38	27:50	1.00	1.084	1.00	1.00	1.084	50.00	50.00	0.439	0.439	1.00
39	32:00	1.00	1.247	1.00	1.00	1.247	50.00	50.00	0.862	0.862	1.00
40	33:17	1.00	1.297	1.00	1.00	1.297	50.00	50.00	0.477	0.477	1.00
41	32:15	1.00	1.256	1.00	1.00	1.256	50.00	50.00	0.559	0.559	1.00
42	30:27	1.00	1.187	1.00	1.00	1.187	50.00	50.00	0.869	0.869	1.00

RIC
04/27/90 8:21:00
SAMPLE: CLP,,,VSTD 50,L,5,22554,U,CC-050,,5ML,,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
RANGE: G 1, 840 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: U3411 #1
CALI: U3411 #2
SCANS 1 TO 840

43161

160101



Quantitation Report File: U3411

04/27/90
(Ref)

Data: U3411.TI

04/27/90 8:21:00

Sample: CLP,,,VSTD 50, L,S,22554,V,CC-050,,5ML,,

Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

Formula: - Instrument: U Weight: 0.011

Submitted by: VERSAR Analyst: MAT Acct. No.: -

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)

Resp. fac. from Library Entry

No	Name
1	C101 BROMOCHLOROMETHANE **IS#1**
2	C010 CHLOROMETHANE
3	C015 BROMOMETHANE
4	C020 VINYL CHLORIDE
5	C025 CHLOROETHANE
6	C030 METHYLENE CHLORIDE
7	C035 ACETONE
8	C040 CARBON DISULFIDE
9	C045 1,1-DICHLOROETHENE
10	C050 1,1-DICHLOROETHANE
11	C053 1,2-DICHLOROETHENE (TOTAL)
12	C060 CHLOROFORM
13	C065 1,2-DICHLOROETHANE
14	CS15 1,2-DICHLOROETHANE-D4 **SS#1**
15	C043 TRICHLOROFLUOROMETHANE
16	CI10 1,4-DIFLUOROBENZENE **IS#2**
17	C110 2-BUTANONE
18	C125 VINYL ACETATE
19	C120 CARBON TETRACHLORIDE
20	C130 BROMODICHLOROMETHANE
21	C140 1,2-DICHLOROPROPANE
22	C145 CIS-1,3-DICHLOROPROPENE
23	C150 TRICHLOROETHENE
24	C165 BENZENE
25	C155 DIBROMOCHLOROMETHANE
26	C160 1,1,2-TRICHLOROETHANE
27	C170 TRANS-1,3-DICHLOROPROPENE
28	C180 BROMOFORM
29	C115 1,1,1-TRICHLOROETHANE
30	CS05 TOLUENE-D8 **SS#2**
31	CI20 CHLOROBENZENE-D5 **IS#3**
32	C205 4-METHYL-2-PENTANONE
33	C210 2-HEXANONE
34	C220 TETRACHLOROETHENE
35	C225 1,1,2,2-TETRACHLOROETHANE
36	C230 TOLUENE
37	C235 CHLOROBENZENE
38	C240 ETHYLBENZENE
39	C245 STYRENE
40	C250 ORTHO & PARA XYLENE
41	C251 META-XYLENE
42	CS10 BROMOFLUOROBENZENE **SS#3**

100102

43411

(Ran)

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	UG/L*	%Tot
1	128	252	10:30	1	1.000	A BB	80565.	50.000	UG/L*	2.38
2	50	52	2:10	1	0.206	A BB	106398.	50.000	UG/L	2.38
3	94	80	3:20	1	0.317	A BB	82997.	50.000	UG/L	2.38
4	62	101	4:12	1	0.401	A BB	85142.	50.000	UG/L	2.38
5	64	126	5:15	1	0.500	A BB	62501.	50.000	UG/L	2.38
6	84	178	7:25	1	0.706	A BB	102191.	50.000	UG/L	2.38
7	43	192	8:00	1	0.762	A BB	39825.	50.000	UG/L	2.38
8	76	216	9:00	1	0.857	A BB	63203.	50.000	UG/L	2.38
9	96	242	10:05	1	0.960	A BB	79942.	50.000	UG/L	2.38
10	63	273	11:22	1	1.083	A BB	195421.	50.000	UG/L	2.38
11	96	289	12:02	1	1.147	A BB	96122.	50.000	UG/L	2.38
12	83	303	12:37	1	1.202	A BB	209318.	50.000	UG/L	2.38
13	62	320	13:20	1	1.270	A BB	170736.	50.000	UG/L	2.38
14	65	318	13:15	1	1.262	A BB	161803.	50.000	UG/L%	2.38
15	101	227	9:27	1	0.901	A BB	162247.	50.000	UG/L	2.38
16	114	497	20:42	16	1.000	A BB	341843.	50.000	UG/L*	2.38
17	72	318	13:15	1	1.262	A BB	15888.	50.000	UG/L	2.38
18	43	363	15:07	16	0.730	A BB	211658.	50.000	UG/L	2.38
19	117	362	15:05	16	0.728	A VB	149107.	50.000	UG/L	2.38
20	83	372	15:30	16	0.748	A BB	200095.	50.000	UG/L	2.38
21	63	406	16:55	16	0.817	A BB	130554.	50.000	UG/L	2.38
22	75	411	17:07	16	0.827	A BB	230589.	50.000	UG/L	2.38
23	130	424	17:40	16	0.853	A BB	159213.	50.000	UG/L	2.38
24	78	437	18:12	16	0.879	A BB	308136.	50.000	UG/L	2.38
25	129	439	18:17	16	0.883	A BB	180282.	50.000	UG/L	2.38
26	97	442	18:25	16	0.889	A BB	116753.	50.000	UG/L	2.38
27	75	441	18:22	16	0.887	A BB	123440.	50.000	UG/L	2.38
28	173	503	20:57	16	1.012	A BB	168586.	50.000	UG/L	2.38
29	97	353	14:42	16	0.710	A BB	169724.	50.000	UG/L	2.38
30	98	587	24:27	31	0.953	A BB	387407.	50.000	UG/L%	2.38
31	117	616	25:40	31	1.000	A BB	337038.	50.000	UG/L*	2.38
32	43	516	21:30	31	0.838	A VV	244267.	50.000	UG/L	2.38
33	43	553	23:02	31	0.898	A BB	195134.	50.000	UG/L	2.38
34	164	559	23:17	31	0.907	A BB	162119.	50.000	UG/L	2.38
35	83	559	23:17	31	0.907	A BB	236925.	50.000	UG/L	2.38
36	92	591	24:37	31	0.959	A BB	214995.	50.000	UG/L	2.38
37	112	619	25:47	31	1.005	A BB	300539.	50.000	UG/L	2.38
38	106	671	27:57	31	1.089	A BB	138636.	50.000	UG/L	2.38
39	104	778	32:25	31	1.263	A BB	259648.	50.000	UG/L	2.38
40	106	811	33:47	31	1.317	A BB	147456.	50.000	UG/L	2.38
41	106	785	32:42	31	1.274	A BB	170247.	50.000	UG/L	2.38
42	95	738	30:45	31	1.198	A BB	253864.	50.000	UG/L%	2.38

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R.Fac	R.Fac(L)	Ratio
1	10:30	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
2	2:10	1.00	0.206	1.00	50.00	50.00	1.321	1.321	1.00
3	3:20	1.00	0.317	1.00	50.00	50.00	1.030	1.030	1.00
4	4:12	1.00	0.401	1.00	50.00	50.00	1.057	1.057	1.00
5	5:15	1.00	0.500	1.00	50.00	50.00	0.776	0.776	1.00
6	7:25	1.00	0.706	1.00	50.00	50.00	1.268	1.268	1.00
7	8:00	1.00	0.762	1.00	50.00	50.00	0.494	0.494	1.00
8	9:00	1.00	0.857	1.00	50.00	50.00	0.784	0.784	1.00
9	10:05	1.00	0.960	1.00	50.00	50.00	0.992	0.992	1.00
10	11:22	1.00	1.083	1.00	50.00	50.00	2.426	2.426	1.00
11	12:02	1.00	1.147	1.00	50.00	50.00	1.193	1.193	1.00
12	12:37	1.00	1.202	1.00	50.00	50.00	2.598	2.598	1.00

100103

U 3411

GlobeWest

(Red)

No	Ret(L)	RRT(L)	Ratio	Amnt	R. Fac	R. Fac(L)	Ratio
13	13; 20	1.00	1.270	1.00	50.00	50.00	1.00
14	13; 15	1.00	1.262	1.00	50.00	50.00	1.00
15	9; 27	1.00	0.901	1.00	50.00	50.00	1.00
16	20; 42	1.00	1.000	1.00	50.00	50.00	1.00
17	13; 15	1.00	1.262	1.00	50.00	50.00	1.00
18	15; 07	1.00	0.730	1.00	50.00	50.00	1.00
19	15; 05	1.00	0.728	1.00	50.00	50.00	1.00
20	15; 30	1.00	0.748	1.00	50.00	50.00	1.00
21	16; 55	1.00	0.817	1.00	50.00	50.00	1.00
22	17; 07	1.00	0.827	1.00	50.00	50.00	1.00
23	17; 40	1.00	0.853	1.00	50.00	50.00	1.00
24	18; 12	1.00	0.879	1.00	50.00	50.00	1.00
25	18; 17	1.00	0.883	1.00	50.00	50.00	1.00
26	18; 25	1.00	0.889	1.00	50.00	50.00	1.00
27	18; 22	1.00	0.887	1.00	50.00	50.00	1.00
28	20; 57	1.00	1.012	1.00	50.00	50.00	1.00
29	14; 42	1.00	0.710	1.00	50.00	50.00	1.00
30	24; 27	1.00	0.953	1.00	50.00	50.00	1.00
31	25; 40	1.00	1.000	1.00	50.00	50.00	1.00
32	21; 30	1.00	0.838	1.00	50.00	50.00	1.00
33	23; 02	1.00	0.898	1.00	50.00	50.00	1.00
34	23; 17	1.00	0.907	1.00	50.00	50.00	1.00
35	23; 17	1.00	0.907	1.00	50.00	50.00	1.00
36	24; 37	1.00	0.959	1.00	50.00	50.00	1.00
37	25; 47	1.00	1.005	1.00	50.00	50.00	1.00
38	27; 57	1.00	1.089	1.00	50.00	50.00	1.00
39	32; 25	1.00	1.263	1.00	50.00	50.00	1.00
40	33; 47	1.00	1.317	1.00	50.00	50.00	1.00
41	32; 42	1.00	1.274	1.00	50.00	50.00	1.00
42	30; 45	1.00	1.198	1.00	50.00	50.00	1.00

100104

8A
VOLATILE INTERNAL STANDARD AREA SUMMARY

OKAYED
(Red)

Lab Name: VERSAR INC.

Contract: _____

La Code: VERSAR Case No.: R3-7

SAS No.: _____ SDG No.: 1

Lab File ID (Standard): U3386

Date Analyzed: 04/26/90

Instrument ID: U

Time Analyzed: 830

Matrix: (soil/water) SOIL Level: (low/med) LOW Column: (pack/cap) PACK

	IS1 (BCM) AREA #	RT	IS2 (DFB) AREA #	RT	IS3 (CBZ) AREA #	RT
12 HOUR STD	55800	10.54	267000	20.75	254000	25.67
UPPER LIMIT	111600		534000		508000	
LOWER LIMIT	27900		133500		127000	
EPA SAMPLE NO.						
01 2	38600	10.54	165000	20.75	146000	25.67
02 3	38200	10.59	164000	20.75	146000	25.67
03 1MS	23600 *	10.54	59000 *	20.75	32300 *	25.67
04 1MSD	19000 *	10.54	46700 *	20.75	24900 *	25.67
5 VBLK87	62300	10.54	275000	20.70	253000	25.62

IS1 (BCM) = Bromochloromethane
 IS2 (DFB) = 1,4-Difluorobenzene
 IS3 (CBZ) = Chlorobenzene

UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

* Column used to flag internal standard area values with an asterisk

8A
VOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: VERSAR INC.

Contract: _____

La Code: VERSAR Case No.: R3-7

SAS No.: _____ SDG No.: 1

Lab File ID (Standard): U3411

Date Analyzed: 04/27/90

Instrument ID: U

Time Analyzed: 821

Matrix: (soil/water) SOIL Level: (low/med) LOW Column: (pack/cap) PACK

	IS1 (BCM) AREA #	RT	IS2 (DFB) AREA #	RT	IS3 (CBZ) AREA #	RT
12 HOUR STD	80600	10.50	342000	20.70	337000	25.67
UPPER LIMIT	161200		684000		674000	
LOWER LIMIT	40300		171000		168500	
EPA SAMPLE NO.						
01 1	23900 *	10.54	50700 *	20.75	25200 *	25.67
02 5	58700	10.54	242000	20.75	201000	25.67
03 6	53200	10.50	226000	20.70	181000	25.62
04 7	49500	10.54	216000	20.75	182000	25.67
5 VBLK12	70100	10.54	276000	21.00	266000	26.04

IS1 (BCM) = Bromochloromethane
 IS2 (DFB) = 1,4-Difluorobenzene
 IS3 (CBZ) = Chlorobenzene

UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

* Column used to flag internal standard area values with an asterisk

Versar[®] Inc.

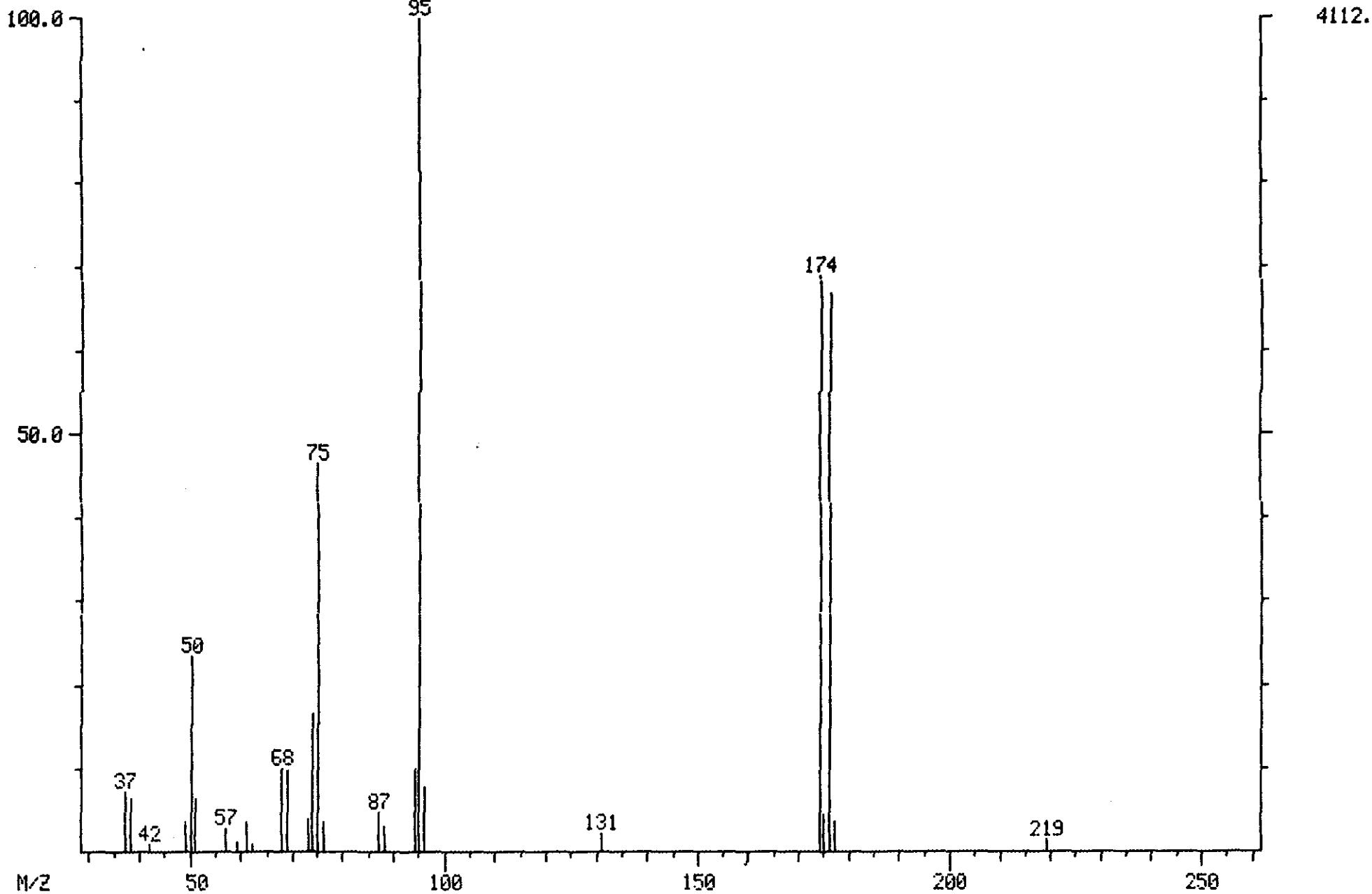
V. RAW QC DATA

100107

MASS SPECTRUM
04/23/90 15:43:00 + 9:12
SAMPLE: VOA BFB 50NG 2UL
COND.: INSTRUMENT U, COLUMN:SP-1000 ISOTHERMO @225C
** NAME: BROMOFLUOROBENZENE
ENHANCED (S 15B 2N 8T)

DATA: U3326 #221
CALI: U3326 #2

BASE M/Z: 95
RIC: 17280.



ORIGINL

Mass

List 04/23/90 15:49:00 + 9:12

Sample: VDA BFB 50NG 2UL

Conds.: INSTRUMENT U, COLUMN: SP-1000 ISOTHERMO @225C

Enhanced (S 15B 2N OT)

Data: U3326 # 221

Cali: U3326 # 2

Base m/z:

9800

RIC: 17280.

Mass	% RA	% RIC	0.	Minima	Min inten:	0.
			#	0 Maxima	Inten.	
37.00?	7.32	1.74		301.		
38.00?	6.25	1.49		257.		
42.00?	0.80	0.19		33.		
47.00?	3.55	0.84		146.		
50.00?	23.42	5.57		963.		
51.00?	6.23	1.48		256.		
57.00?	2.70	0.64		111.		
59.00?	1.00	0.24		41.		
61.00?	3.60	0.86		148.		
62.00?	0.75	0.18		31.		
68.00?	9.97	2.37		410.		
69.00	9.53	2.27		392.		
73.00	3.89	0.93		160.		
74.00	16.51	3.93		679.		
75.00	46.50	11.06		1912.		
76.00	3.60	0.86		148.		
87.00	4.74	1.13		195.		
88.00	3.11	0.74		128.		
94.00	10.00	2.38		411.		
95.00	100.00	23.80		4112.		
96.00	7.86	1.87		323.		
131.00	2.14	0.51		88.		
174.00	68.77	16.37		2828.		
175.00	4.52	1.08		186.		
176.00	66.93	15.93		2752.		
177.00	3.53	0.84		145.		
219.00	1.29	0.31		53.		

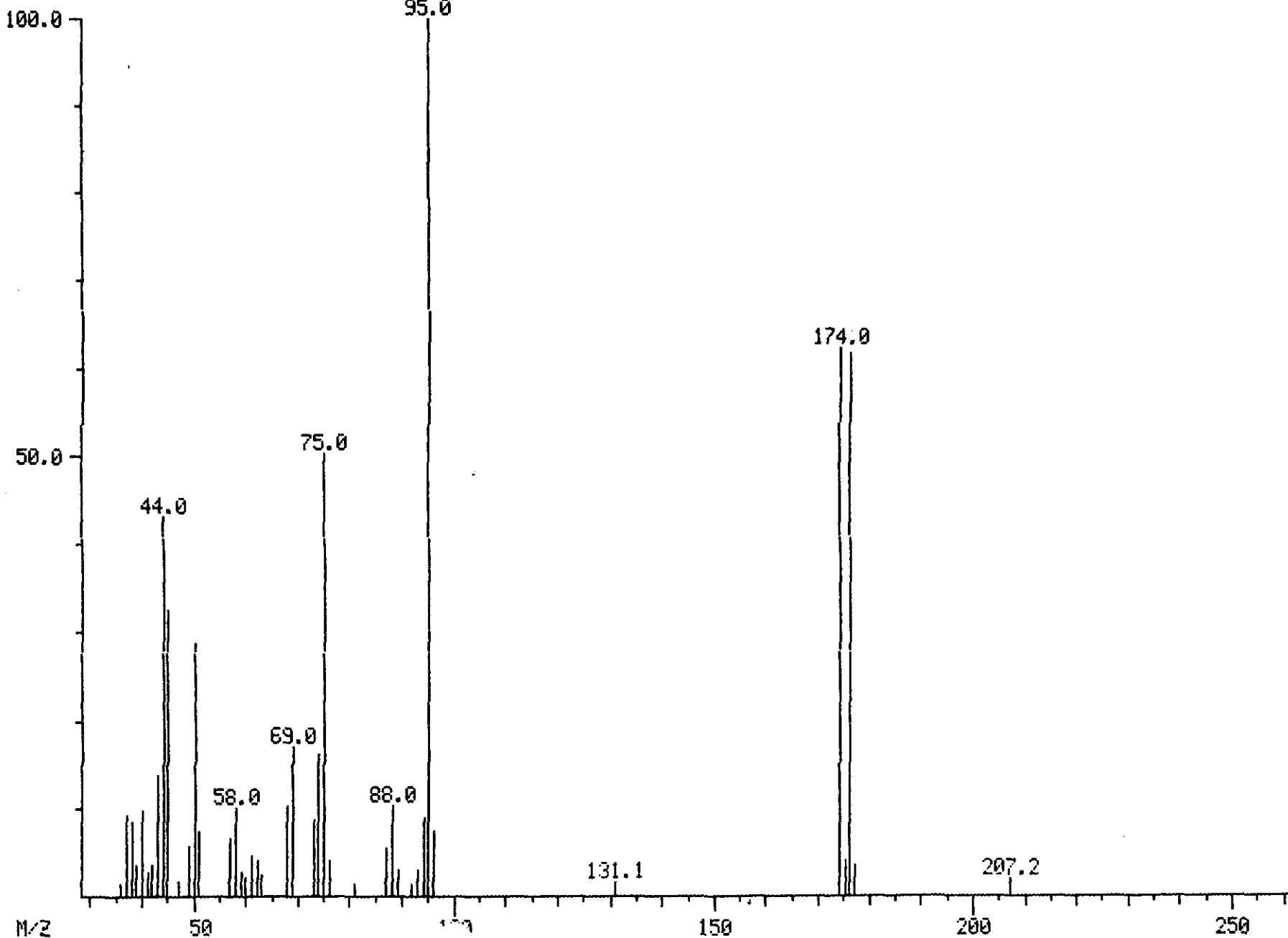
100109

MASS SPECTRUM
04/26/90 7:52:00 + 10:30
SAMPLE: VOA BFB 50NG 2UL
COND.: INSTRUMENT U,COLUMN:SP-1000 ISOTHERMO @225C
** NAME: BROMOFLUOROBENZENE

DATA: U3385 #252
CALI: U3385 #2

BASE M/Z: 95
RIC: 34944.

100 110
5960.



ORIGINAL
(Red)

Mass List
04/26/90 7:52:00 + 10:30
Sample: VOA BFB SONG 2UL
Conds.: INSTRUMENT U, COLUMN: SP-1000 ISOTHERMO @225C
#252

Data: U3385 # 252
Cali: U3385 # 2

Base m/z: 95
RIC: 34944.

Mass	0.00	0.00	O. Minima	Min inten:	O.
	% RA	% RIC	# 0 Maxima	Inten.	
36.00?	1.44	0.25	86.		
37.00?	9.06	1.55	540.		
38.00?	8.37	1.43	499.		
39.00?	3.54	0.60	211.		
40.00?	9.75	1.66	581.		
41.00?	2.75	0.47	164.		
42.00?	3.59	0.61	214.		
43.00?	13.66	2.33	814.		
44.00?	42.95	7.33	2560.		
45.00?	32.55	5.55	1940.		
47.00?	1.71	0.29	102.		
49.00?	5.69	0.97	339.		
50.00?	28.69	4.89	1710.		
51.00?	7.27	1.24	433.		
57.00?	6.48	1.10	386.		
58.00?	9.87	1.68	588.		
59.00?	2.73	0.47	163.		
60.00?	2.18	0.37	130.		
61.00?	4.58	0.78	273.		
62.00?	3.91	0.67	233.		
63.00?	2.55	0.43	152.		
68.00?	10.12	1.73	603.		
69.00	16.90	2.88	1007.		
73.00	8.66	1.48	516.		
74.00	16.17	2.76	964.		
75.00	50.20	8.56	2992.		
76.00	4.06	0.69	242.		
81.00	1.38	0.23	82.		
87.00	5.32	0.91	317.		
88.00	10.15	1.73	605.		
89.00	3.05	0.52	182.		
92.00	1.33	0.23	79.		
93.00	3.09	0.53	184.		
94.00	8.94	1.53	533.		
95.00	100.00	17.06	5960.		
96.00	7.25	1.24	432.		
131.00	1.58	0.27	94.		
174.00	62.21	10.61	3708.		
175.00	4.04	0.69	241.		
176.00	61.61	10.51	3672.		
177.00	3.54	0.60	211.		
207.00	1.85	0.31	110.		

100111

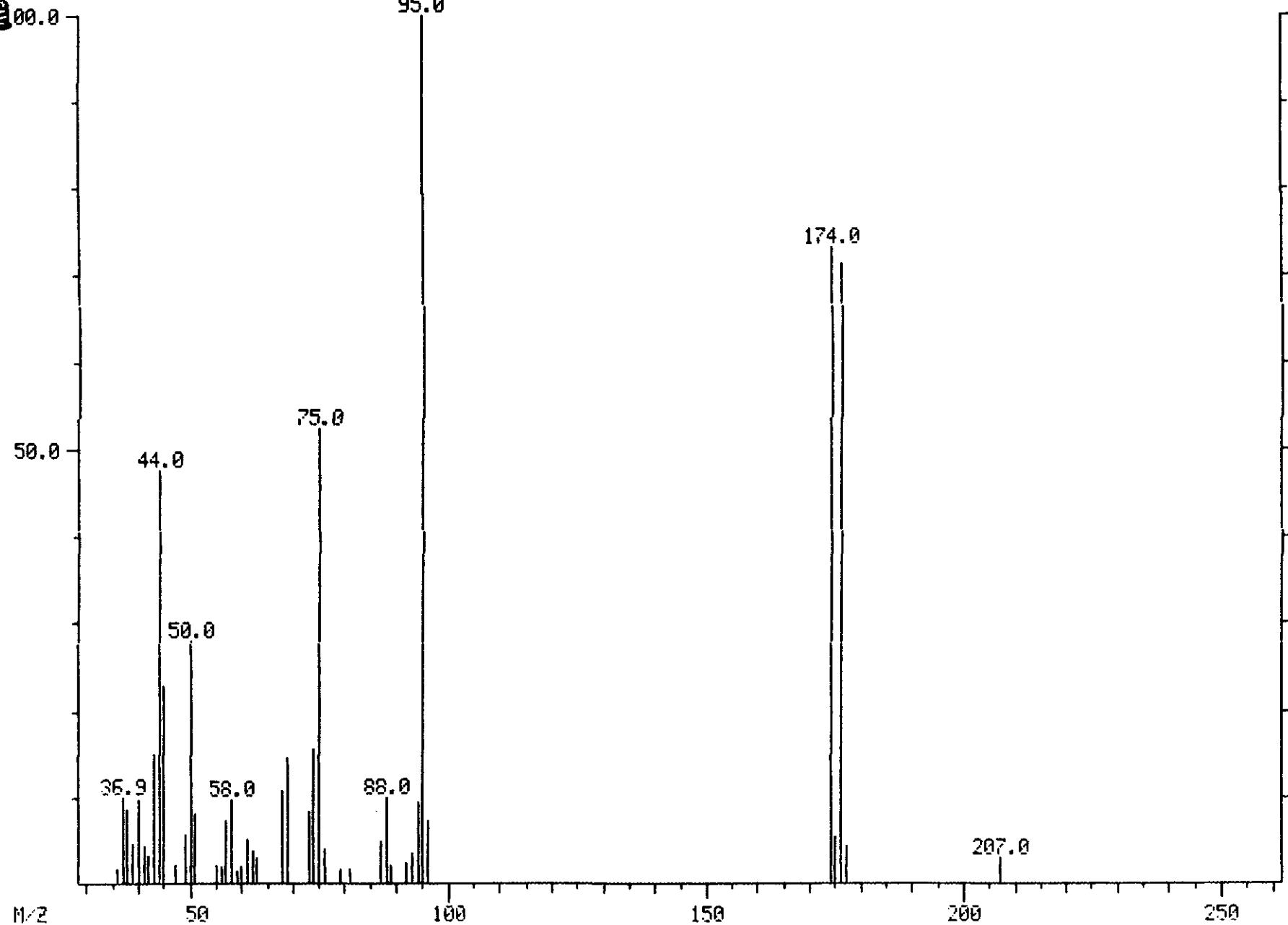
MASS SPECTRUM
04/27/90 7:45:00 + 9:02
SAMPLE: UOA BFB 50NG 2UL

COND.: INSTRUMENT U,COLUMN:SP-1000 ISOTHERMO @225C
** NAME: BROMOFLUOROBENZENE

DATA: U3410 #217
CALI: U3410 #2

BASE M/Z: 95
RIC: 48448.

ORIGINAL
Relab



7886.1 100112

ORIGINAL
(Red)

Mass List

04/27/90 7:45:00 + 9:02

Sample: VOA BFB 50NG 2UL

Conds.: INSTRUMENT U, COLUMN: SP-1000 ISOTHERMO @225C
#217

Data: U3410 # 217

Cali: U3410 # 2

Base m/z: 95

RIC: 48448.

Mass	36	0.00	0.00	O. Minima	Min inten:	O.
	207	% RA	% RIC	# O Maxima	Inten.	
36.00?	1.68	0.27		132.		
37.00?	9.82	1.60		774.		
38.00?	8.71	1.42		686.		
39.00?	4.52	0.73		356.		
40.00?	9.61	1.56		757.		
41.00?	4.40	0.72		347.		
42.00?	3.17	0.52		250.		
43.00?	15.18	2.47		1196.		
44.00?	47.72	7.76		3760.		
45.00?	22.94	3.73		1808.		
47.00?	2.21	0.36		174.		
49.00?	5.67	0.92		447.		
50.00?	28.02	4.56		2208.		
51.00?	8.03	1.31		633.		
55.00?	2.27	0.37		179.		
56.00?	1.88	0.31		148.		
57.00?	7.28	1.18		574.		
58.00?	9.80	1.59		772.		
59.00?	1.33	0.22		105.		
60.00?	1.92	0.31		151.		
61.00?	5.23	0.85		412.		
62.00?	3.77	0.61		297.		
63.00?	2.98	0.49		235.		
68.00?	10.71	1.74		844.		
69.00	14.54	2.37		1146.		
73.00	8.44	1.37		665.		
74.00	15.69	2.55		1236.		
75.00	52.39	8.52		4128.		
76.00	4.05	0.66		319.		
79.00	1.60	0.26		126.		
81.00	1.66	0.27		131.		
87.00	4.77	0.78		376.		
88.00	9.94	1.62		783.		
89.00	2.09	0.34		165.		
92.00	2.39	0.39		188.		
93.00	3.60	0.59		284.		
94.00	9.40	1.53		741.		
95.00	100.00	16.26		7880.		
96.00	7.21	1.17		568.		
174.00	73.10	11.89		5760.		
175.00	5.51	0.90		434.		
176.00	71.37	11.61		5624.		
177.00	4.42	0.72		348.		
207.00	2.86	0.46		225.		

100113

ORIGINAL
(Red)

Versar[®] Inc.

REAGENT BLANK DATA

• 00114

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
VBLK87 (Red)

Lab Name: VERSAR INC.

Contract: _____

Lab Code: VERSAR Case No.: R3-7

SAS No.: _____ SDG No.: 1

Matrix: (soil/water) SOIL

Lab Sample ID: VBLK87

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: U3387

Level: (low/med) LOW

Date Received: _____

% Moisture: not dec. _____

Date Analyzed: 04/26/90

Column: (pack/cap) PACK

Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG	Q
74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene chloride	5	U
67-64-1-----	Acetone	10	U
75-15-0-----	Carbon disulfide	5	U
75-35-4-----	1,1-Dichloroethene	5	U
75-34-3-----	1,1-Dichloroethane	5	U
540-59-0-----	1,2-Dichloroethene (total)	5	U
67-66-3-----	Chloroform	5	U
107-06-2-----	1,2-Dichloroethane	5	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	5	U
56-23-5-----	Carbon tetrachloride	5	U
108-05-4-----	Vinyl acetate	10	U
75-27-4-----	Bromodichloromethane	5	U
78-87-5-----	1,2-Dichloropropane	5	U
10061-01-5-----	cis-1,3-Dichloropropene	5	U
79-01-6-----	Trichloroethene	5	U
124-48-1-----	Dibromochloromethane	5	U
79-00-5-----	1,1,2-Trichloroethane	5	U
71-43-2-----	Benzene	5	U
10061-02-6-----	Trans-1,3-dichloropropene	5	U
75-25-2-----	Bromoform	5	U
108-10-1-----	4-Methyl-2-pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	5	U
79-34-5-----	1,1,2,2-Tetrachloroethane	5	U
108-88-3-----	Toluene	5	U
108-90-7-----	Chlorobenzene	5	U
100-41-4-----	Ethylbenzene	5	U
100-42-5-----	Styrene	5	U
1330-20-7-----	Total xylenes	5	U

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

VBLK87

ORIGINAL
(Red)

Lab Name: VERSAR INC.

Contract: _____

Lab Code: VERSAR Case No.: R3-7 SAS No.: _____ SDG No.: 1 _____

Matrix: (soil/water) SOIL

Lab Sample ID: VBLK87 _____

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: U3387 _____

Level: (low/med) LOW

Date Received: _____

* Moisture: not dec. _____

Date Analyzed: 04/26/90

Column (pack/cap) PACK

Dilution Factor: 1.0 _____

Number TICs found: 0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

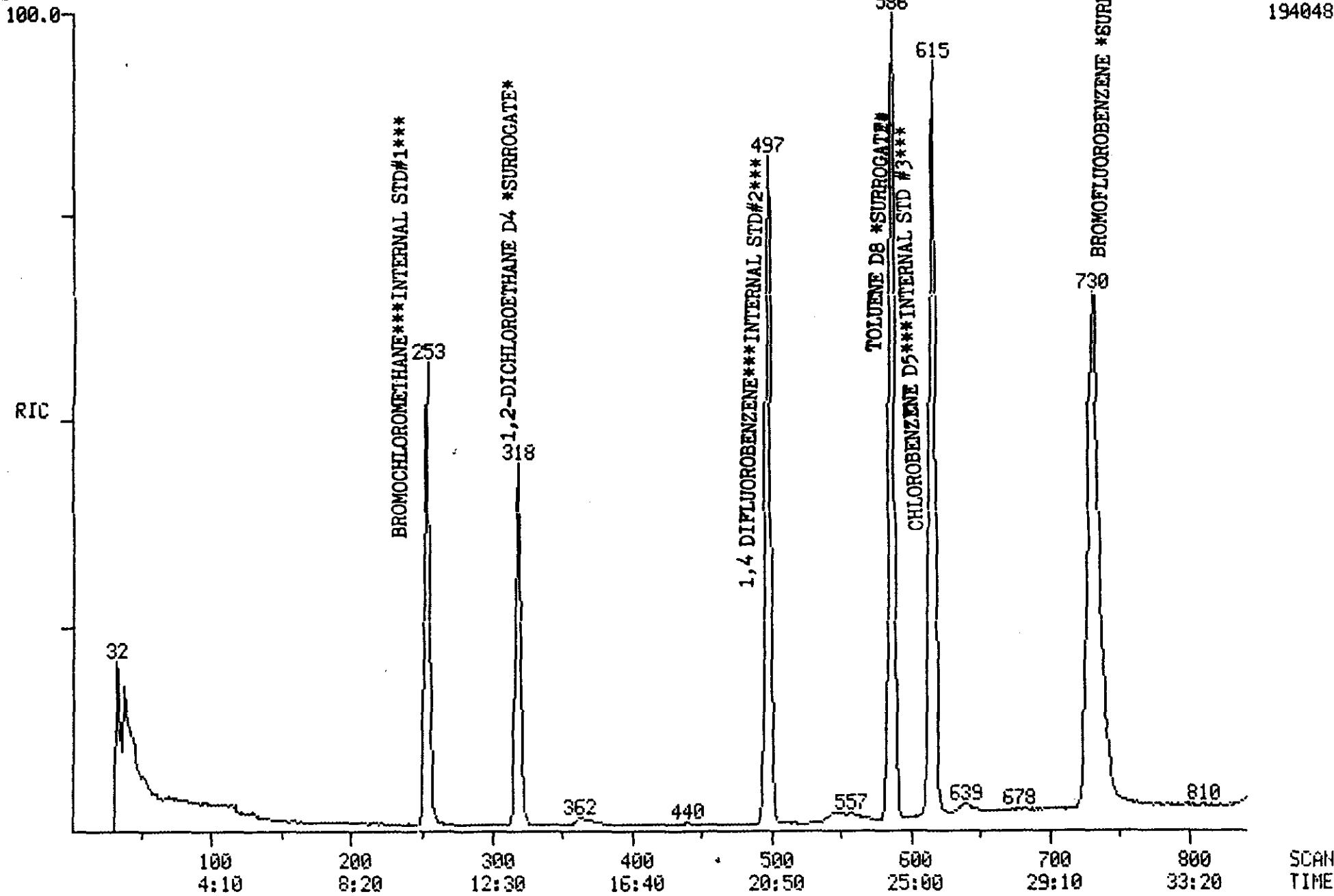
ORIGINALE

RIC
04/26/90 9:31:00
SAMPLE: CLP,,,VBL
COND.: INSTRUMENT
RANGE: G 1, 840

DATA: U3387 #1
CALI: U3387 #2

SCANS 1 TO 840

117



Quantitation Report File: U3387

ORIGINAL
(Red)

Data: U3387.TI

04/26/90 9:31:00

Sample: CLP,,,VBLK87,L,S,VBLK87,V,BLANK,,5 ML.,,

Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

Formula: -

Instrument: U

Weight: 0.011

Submitted by: VERSAR

Analyst: MAT

Acct. No.: -

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)

Resp. fac. from Library Entry

No	Name
1	C101 BROMOCHLOROMETHANE **IS#1**
2	C010 CHLOROMETHANE
3	C015 BROMOMETHANE
4	C020 VINYL CHLORIDE
5	C025 CHLOROETHANE
6	C030 METHYLENE CHLORIDE
7	C035 ACETONE
8	C040 CARBON DISULFIDE
9	C045 1,1-DICHLOROETHENE
10	C050 1,1-DICHLOROETHANE
11	C053 1,2-DICHLOROETHENE (TOTAL)
12	C060 CHLOROFORM
13	C065 1,2-DICHLOROETHANE
14	CS15 1,2-DICHLOROETHANE-D4 **SS#1**
15	C043 TRICHLOROFLUOROMETHANE
16	CI10 1,4-DIFLUOROBENZENE **IS#2**
17	C110 2-BUTANONE
18	C125 VINYL ACETATE
19	C120 CARBON TETRACHLORIDE
20	C130 BROMODICHLOROMETHANE
21	C140 1,2-DICHLOROPROPANE
22	C145 CIS-1,3-DICHLOROPROPENE
23	C150 TRICHLOROETHENE
24	C165 BENZENE
25	C155 DIBROMOCHLOROMETHANE
26	C160 1,1,2-TRICHLOROETHANE
27	C170 TRANS-1,3-DICHLOROPROPENE
28	C180 BROMOFORM
29	C115 1,1,1-TRICHLOROETHANE
30	CS05 TOLUENE-D8 **SS#2**
31	CI20 CHLOROBENZENE-D5 **IS#3**
32	C205 4-METHYL-2-PENTANONE
33	C210 2-HEXANONE
34	C220 TETRACHLOROETHENE
35	C225 1,1,2,2-TETRACHLOROETHANE
36	C230 TOLUENE
37	C235 CHLOROBENZENE
38	C240 ETHYLBENZENE
39	C245 STYRENE
40	C250 ORTHO & PARA XYLENE
41	C251 META-XYLENE
42	CS10 BROMOFLUOROBENZENE **SS#3**

INT STD + SURROG COMPLIANT

No HSL's

No L.S.

Ready For Forms

M.A. Templeton

4/26/90

100118

43387

ORIGINAL
(Red)

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	128	253	10:32	1	1.000	A BB	62297.	50.000 UG/L*	17.29
2	NOT FOUND								
3	NOT FOUND								
4	NOT FOUND								
5	NOT FOUND								
6	NOT FOUND								
7	NOT FOUND								
8	NOT FOUND								
9	NOT FOUND								
10	63	254	10:35	1	1.004	A BB	252.	-0.071 UG/L	BRLO.02
11	NOT FOUND								
12	NOT FOUND								
13	NOT FOUND								
14	65	318	13:15	1	1.257	A BB	124310.	43.612 UG/L%	15.08 87%
15	NOT FOUND								
16	114	497	20:42	16	1.000	A BB	275411.	50.000 UG/L*	17.29
17	NOT FOUND								
18	43	363	15:07	16	0.730	A BB	280.	-0.064 UG/L	BRLO.02
19	NOT FOUND								
20	NOT FOUND								
21	NOT FOUND								
22	NOT FOUND								
23	NOT FOUND								
24	NOT FOUND								
25	NOT FOUND								
26	NOT FOUND								
27	NOT FOUND								
28	NOT FOUND								
29	NOT FOUND								
30	98	586	24:25	31	0.953	A BB	294055.	48.410 UG/L%	16.74 97%
31	117	615	25:37	31	1.000	A BB	252620.	50.000 UG/L*	17.29
32	43	515	21:27	31	0.837	A BB	737.	-0.158 UG/L	BRLO.05
33	43	553	23:02	31	0.899	A BB	359.	-0.094 UG/L	BRLO.03
34	NOT FOUND								
35	NOT FOUND								
36	92	590	24:35	31	0.959	A BB	666.	0.192 UG/L	BRLO.07
37	NOT FOUND								
38	NOT FOUND								
39	NOT FOUND								
40	NOT FOUND								
41	NOT FOUND								
42	95	730	30:25	31	1.187	A BB	204931.	46.677 UG/L%	16.14 93%

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
1	10:32	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
2	2:12		0.209						
3	3:22		0.320						
4	4:15		0.403						
5	5:15		0.498						
6	7:27		0.708						
7	8:02		0.763						
8	9:02		0.858						
9	10:07		0.960						
10	11:25	0.93	1.083	0.93	0.07	50.00	0.004	2.837	0.00
11	12:05		1.146						
12	12:40		1.202						

100119

43387

ORIGINAL
(Red)

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
13	13:22		1.269						
14	13:17	1.00	1.261	1.00	43.61	50.00	1.995	2.288	0.87
15	9:30		0.901						
16	20:45	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
17	13:17		1.261						
18	15:10	1.00	0.731	1.00	0.06	50.00	0.001	0.791	0.00
19	15:07		0.729						
20	15:32		0.749						
21	16:57		0.817						
22	17:10		0.827						
23	17:42		0.853						
24	18:12		0.878						
25	18:20		0.884						
26	18:25		0.888						
27	18:25		0.888						
28	21:00		1.012						
29	14:42		0.709						
30	24:27	1.00	0.953	1.00	48.41	50.00	1.164	1.202	0.97
31	25:40	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
32	21:30	1.00	0.838	1.00	0.16	50.00	0.003	0.922	0.00
33	23:02	1.00	0.898	1.00	0.09	50.00	0.001	0.759	0.00
34	23:17		0.907						
35	23:17		0.907						
36	24:37	1.00	0.959	1.00	0.19	50.00	0.003	0.686	0.00
37	25:47		1.005						
38	27:50		1.084						
39	32:00		1.247						
40	33:17		1.297						
41	32:15		1.256						
42	30:27	1.00	1.187	1.00	46.68	50.00	0.811	0.869	0.93

100120

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
VBLK12 (Rev)

Lab Name: VERSAR INC.

Contract: _____

Lab Code: VERSAR Case No.: R3-7

SAS No.: _____ SDG No.: 1

Matrix: (soil/water) SOIL

Lab Sample ID: VBLK12

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: U3412

Level: (low/med) LOW

Date Received: _____

* Moisture: not dec. _____

Date Analyzed: 04/27/90

Column: (pack/cap) PACK

Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene chloride	5	U
67-64-1-----	Acetone	10	U
75-15-0-----	Carbon disulfide	5	U
75-35-4-----	1,1-Dichloroethene	5	U
75-34-3-----	1,1-Dichloroethane	5	U
540-59-0-----	1,2-Dichloroethene (total)	5	U
67-66-3-----	Chloroform	5	U
107-06-2-----	1,2-Dichloroethane	5	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	5	U
56-23-5-----	Carbon tetrachloride	5	U
108-05-4-----	Vinyl acetate	10	U
75-27-4-----	Bromodichloromethane	5	U
78-87-5-----	1,2-Dichloropropane	5	U
10061-01-5-----	cis-1,3-Dichloropropene	5	U
79-01-6-----	Trichloroethene	5	U
124-48-1-----	Dibromochloromethane	5	U
79-00-5-----	1,1,2-Trichloroethane	5	U
71-43-2-----	Benzene	5	U
10061-02-6-----	Trans-1,3-dichloropropene	5	U
75-25-2-----	Bromoform	5	U
108-10-1-----	4-Methyl-2-pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	5	U
79-34-5-----	1,1,2,2-Tetrachloroethane	5	U
108-88-3-----	Toluene	5	U
108-90-7-----	Chlorobenzene	5	U
100-41-4-----	Ethylbenzene	5	U
100-42-5-----	Styrene	5	U
1330-20-7-----	Total xylenes	5	U

1E
 VOLATILE ORGANICS ANALYSIS DATA SHEET
 TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ORIGINAL
VBLK12 (Red)

Lab Name: VERSAR INC.

Contract: _____

Lab Code: VERSAR Case No.: R3-7

SAS No.: _____ SDG No.: 1 _____

Matrix: (soil/water) SOIL

Lab Sample ID: VBLK12 _____

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: U3412 _____

Level: (low/med) LOW

Date Received: _____

* Moisture: not dec. _____

Date Analyzed: 04/27/90

Column (pack/cap) PACK

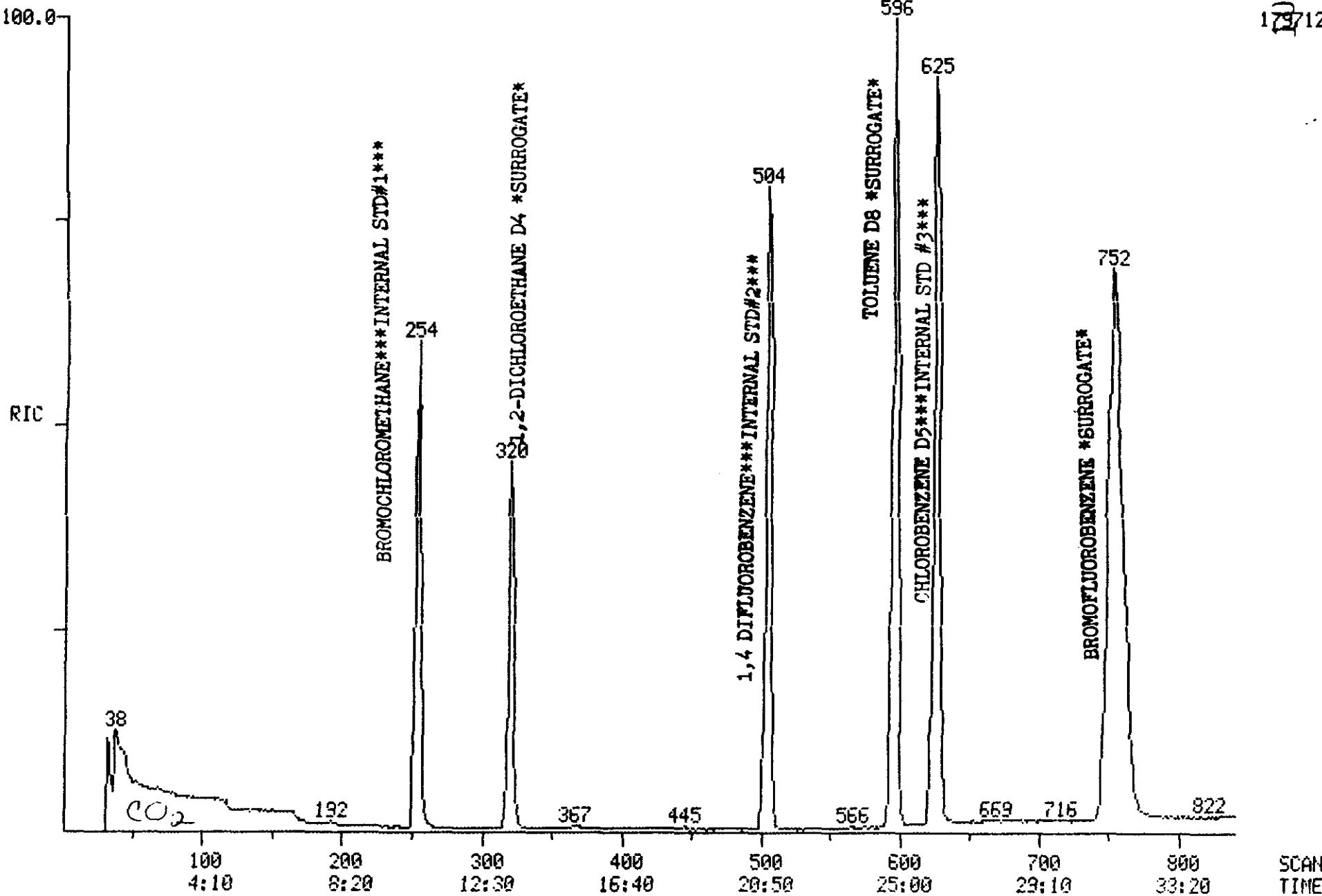
Dilution Factor: 1.0 _____

Number TICs found: 0

CONCENTRATION UNITS:
 (ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

RIC 04/27/90 9:26:00 DHTH: U3412 #1 5LHN# 1 10 840
CALI: U3412 #2 SAMPLE: CLP,,,UBLK12,L,S,UBLK12,U,BLANK,,5 ML,,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
RANGE: G 1, 840 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20,



1212.
N0123

Data: U3412.TI

04/27/90 9:26:00

Sample: CLP,,, VBLK12, L, S, VBLK12, V, BLANK,, 5 ML.,,

Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

Formula: -

Instrument: U

Weight: 0.011

Submitted by: VERSAR

Analyst: MAT

Acct. No.: -

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)

Resp. fac. from Library Entry

No	Name
1	C101 BROMOCHLOROMETHANE **IS#1**
2	C010 CHLOROMETHANE
3	C015 BROMOMETHANE
4	C020 VINYL CHLORIDE
5	C025 CHLOROETHANE
6	C030 METHYLENE CHLORIDE
7	C035 ACETONE
8	C040 CARBON DISULFIDE
9	C045 1,1-DICHLOROETHENE
10	C050 1,1-DICHLOROETHANE
11	C053 1,2-DICHLOROETHENE (TOTAL)
12	C060 CHLOROFORM
13	C065 1,2-DICHLOROETHANE
14	CS15 1,2-DICHLOROETHANE-D4 **SS#1**
15	C043 TRICHLOROFUOROMETHANE
16	CI10 1,4-DIFLUOROBENZENE **IS#2**
17	C110 2-BUTANONE
18	C125 VINYL ACETATE
19	C120 CARBON TETRACHLORIDE
20	C130 BROMODICHLOROMETHANE
21	C140 1,2-DICHLOROPROPANE
22	C145 CIS-1,3-DICHLOROPROPENE
23	C150 TRICHLOROETHENE
24	C165 BENZENE
25	C155 DIBROMOCHLOROMETHANE
26	C160 1,1,2-TRICHLOROETHANE
27	C170 TRANS-1,3-DICHLOROPROPENE
28	C180 BROMOFORM
29	C115 1,1,1-TRICHLOROETHANE
30	CS05 TOLUENE-D8 **SS#2**
31	CI20 CHLOROBENZENE-D5 **IS#3**
32	C205 4-METHYL-2-PENTANONE
33	C210 2-HEXANONE
34	C220 TETRACHLOROETHENE
35	C225 1,1,2,2-TETRACHLOROETHANE
36	C230 TOLUENE
37	C235 CHLOROBENZENE
38	C240 ETHYLBENZENE
39	C245 STYRENE
40	C250 ORTHO & PARA XYLENE
41	C251 META-XYLENE
42	CS10 BROMOFLUOROBENZENE **SS#3**

INT STD + SURRG COMPLIANT

No HSL's

No LS.

Ready For Forms

M. A. Templeton

4/27/90

100124

ORIGINAL
(Red)

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	128	253	10:32	1	1.000	A BB	70104.✓	50.000 UG/L*	17.03
2	NOT FOUND								
3	NOT FOUND								
4	NOT FOUND								
5	NOT FOUND								
6	NOT FOUND								
7	NOT FOUND								
8	NOT FOUND								
9	NOT FOUND								
10	63	254	10:35	1	1.004	A BB	269.	0.079 UG/L	BRLO. 03
11	NOT FOUND								
12	NOT FOUND								
13	NOT FOUND								
14	65	320	13:20	1	1.265	A BB	118849.	42.207 UG/L%	14.37 84
15	NOT FOUND								
16	114	504	21:00	16	1.000	A BB	276334.✓	50.000 UG/L*	17.03
17	NOT FOUND								
18	NOT FOUND								
19	NOT FOUND								
20	NOT FOUND								
21	NOT FOUND								
22	NOT FOUND								
23	NOT FOUND								
24	NOT FOUND								
25	NOT FOUND								
26	NOT FOUND								
27	NOT FOUND								
28	NOT FOUND								
29	NOT FOUND								
30	98	595	24:47	31	0.952	A BB	311994.	51.095 UG/L%	17.40 10.
31	117	625	26:02	31	1.000	A BB	265615.✓	50.000 UG/L*	17.03
32	NOT FOUND								
33	NOT FOUND								
34	NOT FOUND								
35	NOT FOUND								
36	92	600	25:00	31	0.960	A BB	538.	0.159 UG/L	BRLO. 05
37	NOT FOUND								
38	NOT FOUND								
39	NOT FOUND								
40	NOT FOUND								
41	NOT FOUND								
42	95	752	31:20	31	1.203	A BB	200557.	50.122 UG/L%	17.07 100

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
1	10:30	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
2	2:10		0.206						
3	3:20		0.317						
4	4:12		0.401						
5	5:15		0.500						
6	7:25		0.706						
7	8:00		0.762						
8	9:00		0.857						
9	10:05		0.960						
10	11:22	0.93	1.083	0.93	0.08	50.00	0.004	2.426	0.00
11	12:02		1.147						
12	12:37		1.202						

100125

**OPTIONAL
(Red)**

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
13	13:20		1. 270						
14	13:15	1. 01	1. 262	1. 00	42. 21	50. 00	1. 695	2. 008	0. 84
15	9:27		0. 901						
16	20:42	1. 01	1. 000	1. 00	50. 00	50. 00	1. 000	1. 000	1. 00
17	13:15		1. 262						
18	15:07		0. 730						
19	15:05		0. 728						
20	15:30		0. 748						
21	16:55		0. 817						
22	17:07		0. 827						
23	17:40		0. 853						
24	18:12		0. 879						
25	18:17		0. 883						
26	18:25		0. 889						
27	18:22		0. 887						
28	20:57		1. 012						
29	14:42		0. 710						
30	24:27	1. 01	0. 953	1. 00	51. 09	50. 00	1. 175	1. 149	1. 02
31	25:40	1. 01	1. 000	1. 00	50. 00	50. 00	1. 000	1. 000	1. 00
32	21:30		0. 838						
33	23:02		0. 898						
34	23:17		0. 907						
35	23:17		0. 907						
36	24:37	1. 02	0. 959	1. 00	0. 16	50. 00	0. 002	0. 638	0. 00
37	25:47		1. 005						
38	27:57		1. 089						
39	32:25		1. 263						
40	33:47		1. 317						
41	32:42		1. 274						
42	30:45	1. 02	1. 198	1. 00	50. 12	50. 00	0. 755	0. 753	1. 00

100126

Versar^{inc.}

**ORIGINAL
(Red)**

MATRIX SPIKE DATA

100127

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: VERSAR INC.

Contract: _____

1MS	ORIGINAL (Red)
-----	-------------------

Lab Code: VERSAR Case No.: R3-7 SAS No.: _____ SDG No.: 1

Matrix: (soil/water) SOIL Lab Sample ID: 16426MS

Sample wt/vol: 5.0 (g/mL) G Lab File ID: U3398

Level: (low/med) LOW Date Received: 04/19/90

% Moisture: not dec. 30 Date Analyzed: 04/26/90

Column: (pack/cap) PACK Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG	Q
---------	----------	---	---

74-87-3-----	Chloromethane	14	U
74-83-9-----	Bromomethane	14	U
75-01-4-----	Vinyl chloride	14	U
75-00-3-----	Chloroethane	14	U
75-09-2-----	Methylene chloride	11	
67-64-1-----	Acetone	51	
75-15-0-----	Carbon disulfide	7	U
75-35-4-----	1,1-Dichloroethene	7	U
75-34-3-----	1,1-Dichloroethane	7	U
540-59-0-----	1,2-Dichloroethene (total)	7	U
67-66-3-----	Chloroform	7	U
107-06-2-----	1,2-Dichloroethane	7	U
78-93-3-----	2-Butanone	14	U
71-55-6-----	1,1,1-Trichloroethane	7	U
56-23-5-----	Carbon tetrachloride	7	U
108-05-4-----	Vinyl acetate	14	U
75-27-4-----	Bromodichloromethane	7	U
78-87-5-----	1,2-Dichloropropane	7	U
10061-01-5-----	cis-1,3-Dichloropropene	7	U
79-01-6-----	Trichloroethene	7	U
124-48-1-----	Dibromochloromethane	7	U
79-00-5-----	1,1,2-Trichloroethane	7	U
71-43-2-----	Benzene	7	U
10061-02-6-----	Trans-1,3-dichloropropene	7	U
75-25-2-----	Bromoform	7	U
108-10-1-----	4-Methyl-2-pentanone	14	U
591-78-6-----	2-Hexanone	14	U
127-18-4-----	Tetrachloroethene	7	U
79-34-5-----	1,1,2,2-Tetrachloroethane	7	U
108-88-3-----	Toluene	7	U
108-90-7-----	Chlorobenzene	7	U
100-41-4-----	Ethylbenzene	7	U
100-42-5-----	Styrene	7	U
1330-20-7-----	Total xylenes	7	U

OPTIONAL
(Ref)

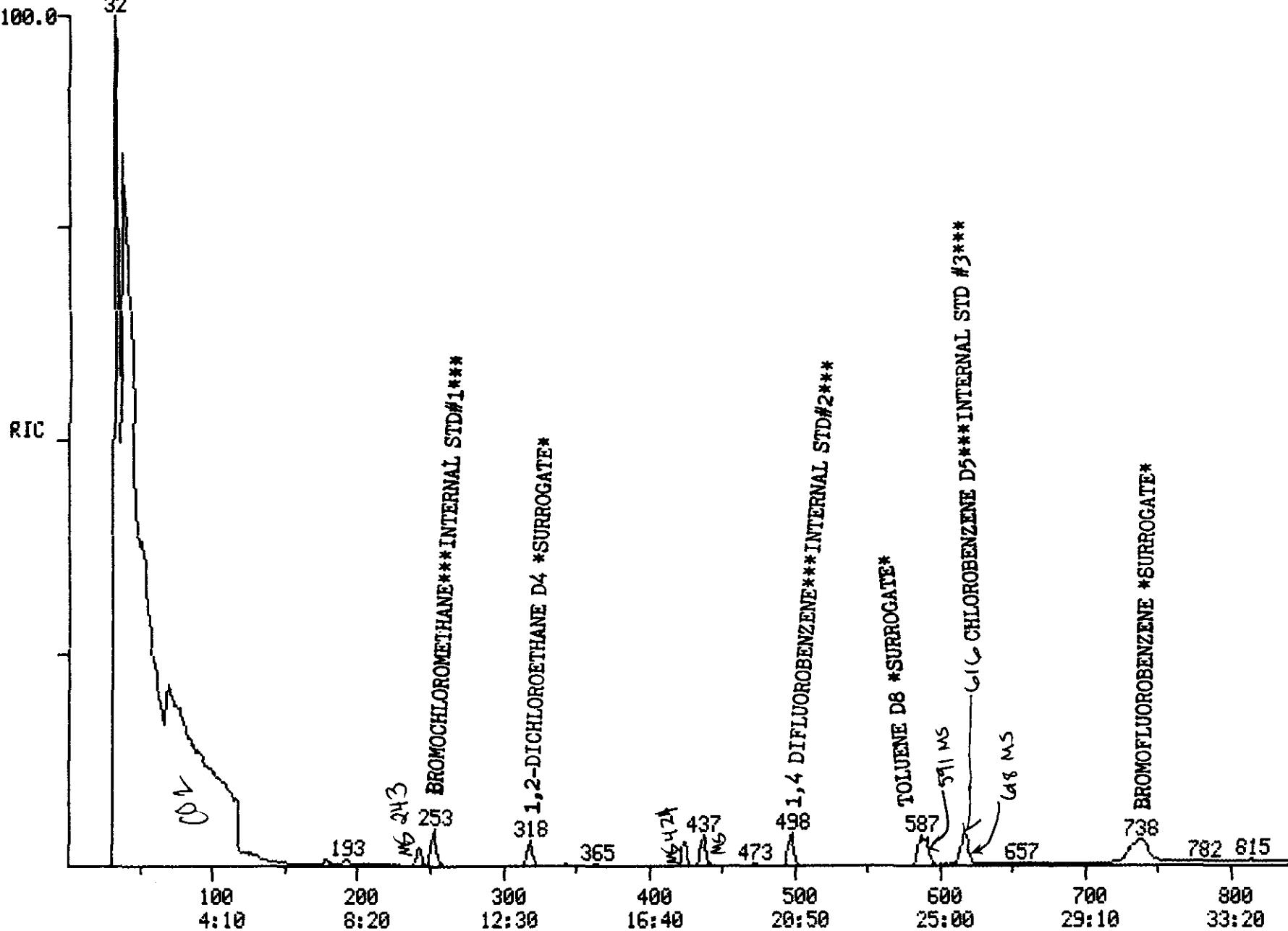
RIC
04/26/90 18:11:00

DATA: U3398 #1
CALI: U3398 #2

SCANS 1 TO 840

SAMPLE: CLP, VERSCOM, 2536, 1MS, L, S, 16426MS, U, , 420.1.0, 2,5G/5ML,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
RANGE: G 1, 840 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

747520.
747520.



OPTIONAL
(Ref)

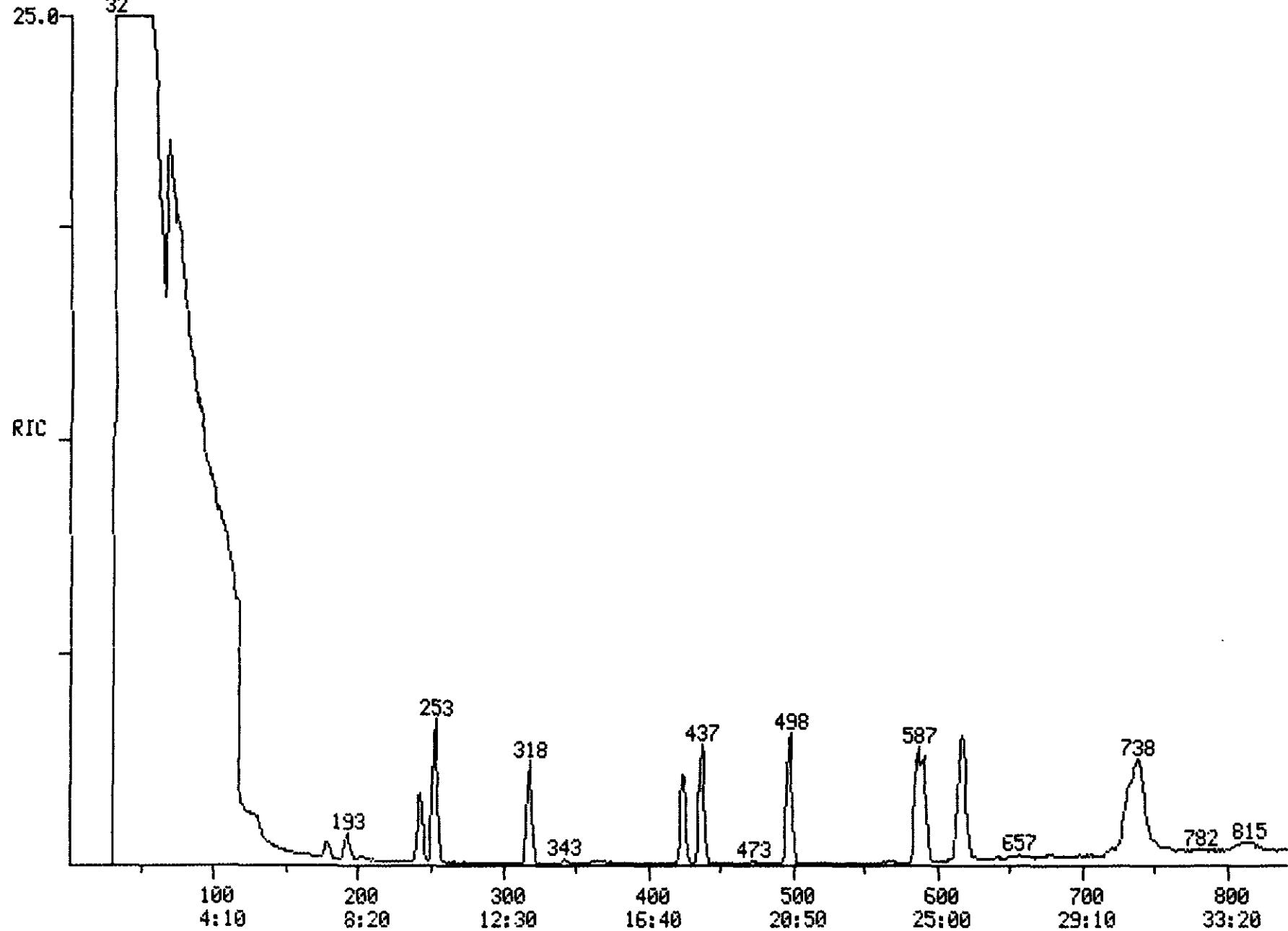
RIC
04/26/90 18:11:00

DATA: U3398 #1
CALI: U3398 #2

SCANS 1 TO 840

SAMPLE: CLP, VERSCOM, 2536, 1MS,L,S,16426MS,U,,420.1.0,2,5G/5ML,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
RANGE: G 1, 840 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

1100130
186880.



Data: U3398.TI

04/26/90 18:11:00

Sample: CLP, VERSCDM, 2536, 1MS, L, S, 16426MS, V., 420.1.0, 2, 5G/5ML,

Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

Formula: ALS 8

Instrument: U

Weight: 0.011

Submitted by: VERSAR

Analyst: DD

Acct. No.: 420.1.0

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)

30.0%

Resp. fac. from Library Entry

No	Name
1	C101 BROMOCHLOROMETHANE **IS#1**
2	C010 CHLOROMETHANE
3	C015 BROMOMETHANE
4	C020 VINYL CHLORIDE
5	C025 CHLOROETHANE
6	C030 METHYLENE CHLORIDE
7	C035 ACETONE
8	C040 CARBON DISULFIDE
9	C045 1,1-DICHLOROETHENE
10	C050 1,1-DICHLOROETHANE
11	C053 1,2-DICHLOROETHENE (TOTAL)
12	C060 CHLOROFORM
13	C065 1,2-DICHLOROETHANE
14	CS15 1,2-DICHLOROETHANE-D4 **SS#1**
15	C043 TRICHLOROFUOROMETHANE
16	CI10 1,4-DIFLUOROBENZENE **IS#2**
17	C110 2-BUTANONE
18	C125 VINYL ACETATE
19	C120 CARBON TETRACHLORIDE
20	C130 BROMODICHLOROMETHANE
21	C140 1,2-DICHLOROPROPANE
22	C145 CIS-1,3-DICHLOROPROPENE
23	C150 TRICHLOROETHENE
24	C165 BENZENE
25	C155 DIBROMOCHLOROMETHANE
26	C160 1,1,2-TRICHLOROETHANE
27	C170 TRANS-1,3-DICHLOROPROPENE
28	C180 BROMOFORM
29	C115 1,1,1-TRICHLOROETHANE
30	CS05 TOLUENE-D8 **SS#2**
31	CI20 CHLOROBENZENE-D5 **IS#3**
32	C205 4-METHYL-2-PENTANONE
33	C210 2-HEXANONE
34	C220 TETRACHLOROETHENE
35	C225 1,1,2,2-TETRACHLOROETHANE
36	C230 TOLUENE
37	C235 CHLOROBENZENE
38	C240 ETHYLBENZENE
39	C245 STYRENE
40	C250 ORTHO & PARA XYLENE
41	C251 META-XYLENE
42	CS10 BROMOFLUOROBENZENE **SS#3**

Ready for

100131

613398

ORIGIN

(Red)

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	(Red)	%Tot	
1	128	253	10:32	1	1.000	A BB	23631.	50.000	UG/L*	9.19	
2	NOT FOUND						15 Nov				
3	NOT FOUND										
4	NOT FOUND										
5	NOT FOUND										
6	84	179✓	7:27	1	0.708	A BB	5099.	7.864	UG/L	1.45	
7	43	192✓	8:00	1	0.759	A BB	12258.	35.683	UG/L	6.56	
8	NOT FOUND										
9	96	243	10:07	1	0.960	A BB	15367.	MS	28.232	UG/L	5.19
10	NOT FOUND										
11	NOT FOUND										
12	NOT FOUND										
13	NOT FOUND										
14	65	318	13:15	1	1.257	A BB	33483.	30.968	UG/L%	5.69	
15	NOT FOUND										
16	114	498	20:45	16	1.000	A BB	59035.	MS	50.000	UG/L*	9.19
17	NOT FOUND						low				
18	NOT FOUND										
19	NOT FOUND										
20	NOT FOUND										
21	NOT FOUND										
22	NOT FOUND										
23	130	424	17:40	16	0.851	A BB	18288.	MS	35.788	UG/L	6.58
24	78	437	18:12	16	0.878	A BB	57749.	MS	54.989	UG/L	10.11
25	NOT FOUND										
26	NOT FOUND										
27	NOT FOUND										
28	NOT FOUND										
29	NOT FOUND										
30	98	587	24:27	31	0.953	A BB	45277.	58.377	UG/L%	10.73	
31	117	616	25:40	31	1.000	A BB	32256.	MS	50.000	UG/L*	9.19
32	NOT FOUND						low				
33	NOT FOUND										
34	NOT FOUND										
35	NOT FOUND										
36	92	591	24:37	31	0.959	A BB	25531.	MS	57.662	UG/L	10.60
37	112	618	25:45	31	1.003	A BB	27677.	MS	46.907	UG/L	8.62
38	NOT FOUND										
39	NOT FOUND										
40	NOT FOUND										
41	NOT FOUND										
42	95	732	30:30	31	1.188	A BB	21041.	MS	37.534	UG/L%	6.90

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
1	10:32	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
2	2:12		0.209						
3	3:22		0.320						
4	4:15		0.403						
5	5:15		0.498						
6	7:27	1.00	0.708	1.00	7.86	50.00	0.216	1.372	0.16
7	8:02	0.99	0.763	0.99	35.68	50.00	0.519	0.727	0.71
8	9:02		0.858						
9	10:07	1.00	0.960	1.00	28.23	50.00	0.650	1.152	0.56
10	11:25		1.083						
11	12:05		1.146						
12	12:40		1.202						

100132

43398
Original
(Red)

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
13	13:22		1. 269						
14	13:17	1. 00	1. 261	1. 00	30. 97	50. 00	1. 417	2. 288	0. 62
15	9:30		0. 901						
16	20:45	1. 00	1. 000	1. 00	50. 00	50. 00	1. 000	1. 000	1. 00
17	13:17		1. 261						
18	15:10		0. 731						
19	15:07		0. 729						
20	15:32		0. 749						
21	16:57		0. 817						
22	17:10		0. 827						
23	17:42	1. 00	0. 853	1. 00	35. 79	50. 00	0. 310	0. 433	0. 72
24	18:12	1. 00	0. 878	1. 00	54. 99	50. 00	0. 978	0. 889	1. 10
25	18:20		0. 884						
26	18:25		0. 888						
27	18:25		0. 888						
28	21:00		1. 012						
29	14:42		0. 709						
30	24:27	1. 00	0. 953	1. 00	58. 38	50. 00	1. 404	1. 202	1. 17
31	25:40	1. 00	1. 000	1. 00	50. 00	50. 00	1. 000	1. 000	1. 00
32	21:30		0. 838						
33	23:02		0. 898						
34	23:17		0. 907						
35	23:17		0. 907						
36	24:37	1. 00	0. 959	1. 00	57. 66	50. 00	0. 792	0. 686	1. 15
37	25:47	1. 00	1. 005	1. 00	46. 91	50. 00	0. 858	0. 915	0. 94
38	27:50		1. 084						
39	32:00		1. 247						
40	33:17		1. 297						
41	32:15		1. 256						
42	30:27	1. 00	1. 187	1. 00	37. 53	50. 00	0. 652	0. 869	0. 75

100133

Versar^{inc.}

ORIGINAL
(Red)

MATRIX SPIKE DUPLICATE DATA

100184

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: VERSAR INC.

Contract: _____

1MSD ORIGINAL
(Red)

Lab Code: VERSAR Case No.: R3-7 SAS No.: _____ SDG No.: 1 _____

Matrix: (soil/water) SOIL

Lab Sample ID: 16426MSD

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: U3399

Level: (low/med) LOW

Date Received: 04/19/90

% Moisture: not dec. 30

Date Analyzed: 04/26/90

Column: (pack/cap) PACK

Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG	Q
74-87-3-----	Chloromethane	14	U
74-83-9-----	Bromomethane	14	U
75-01-4-----	Vinyl chloride	14	U
75-00-3-----	Chloroethane	14	U
75-09-2-----	Methylene chloride	12	
67-64-1-----	Acetone	49	
75-15-0-----	Carbon disulfide	7	U
75-35-4-----	1,1-Dichloroethene	7	U
75-34-3-----	1,1-Dichloroethane	7	U
540-59-0-----	1,2-Dichloroethene (total)	7	U
67-66-3-----	Chloroform	7	U
107-06-2-----	1,2-Dichloroethane	7	U
78-93-3-----	2-Butanone	14	U
71-55-6-----	1,1,1-Trichloroethane	7	U
56-23-5-----	Carbon tetrachloride	7	U
108-05-4-----	Vinyl acetate	14	U
75-27-4-----	Bromodichloromethane	7	U
78-87-5-----	1,2-Dichloropropane	7	U
10061-01-5-----	cis-1,3-Dichloropropene	7	U
79-01-6-----	Trichloroethene	7	U
124-48-1-----	Dibromochloromethane	7	U
79-00-5-----	1,1,2-Trichloroethane	7	U
71-43-2-----	Benzene	7	U
10061-02-6-----	Trans-1,3-dichloropropene	7	U
75-25-2-----	Bromoform	7	U
108-10-1-----	4-Methyl-2-pentanone	14	U
591-78-6-----	2-Hexanone	14	U
127-18-4-----	Tetrachloroethene	7	U
79-34-5-----	1,1,2,2-Tetrachloroethane	7	U
108-88-3-----	Toluene	7	U
108-90-7-----	Chlorobenzene	7	U
100-41-4-----	Ethylbenzene	7	U
100-42-5-----	Styrene	7	U
1330-20-7-----	Total xylenes	7	U

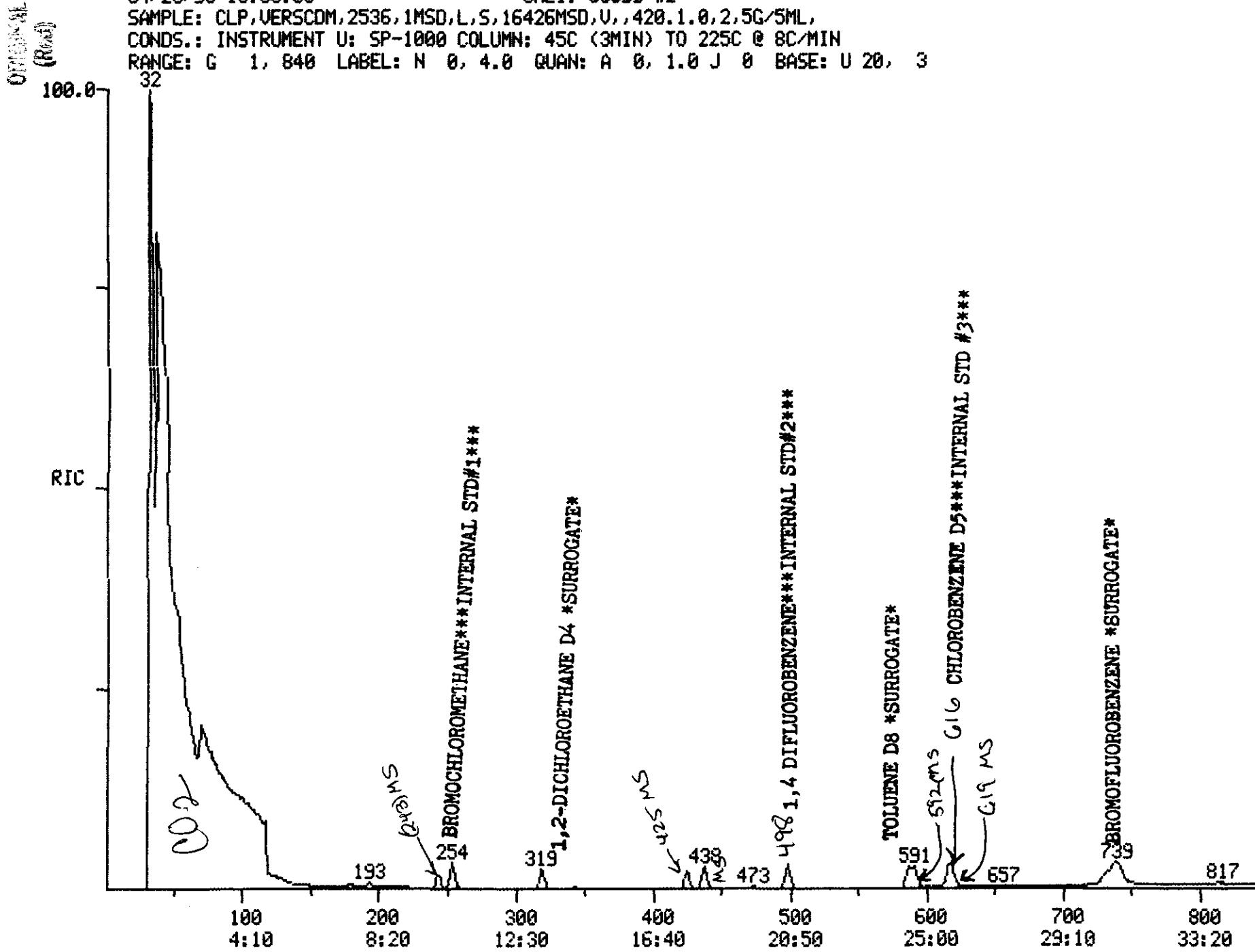
RIC
04/26/90 18:53:00

DATA: U3399 #1
CALI: U3399 #2

SCANS 1 TO 840

SAMPLE: CLP, VERSCDM, 2536, 1MSD,L,S,16426MSD,U,,420.1.0,2,5G/5ML,
COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN
RANGE: G 1, 840 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

757760.
757760.100136



ORIGINAL
(Raw)

RIC
04/26/90 18:53:00

DATA: U3399 #1

SCANS 1 TO 840

CALI: U3399 #2

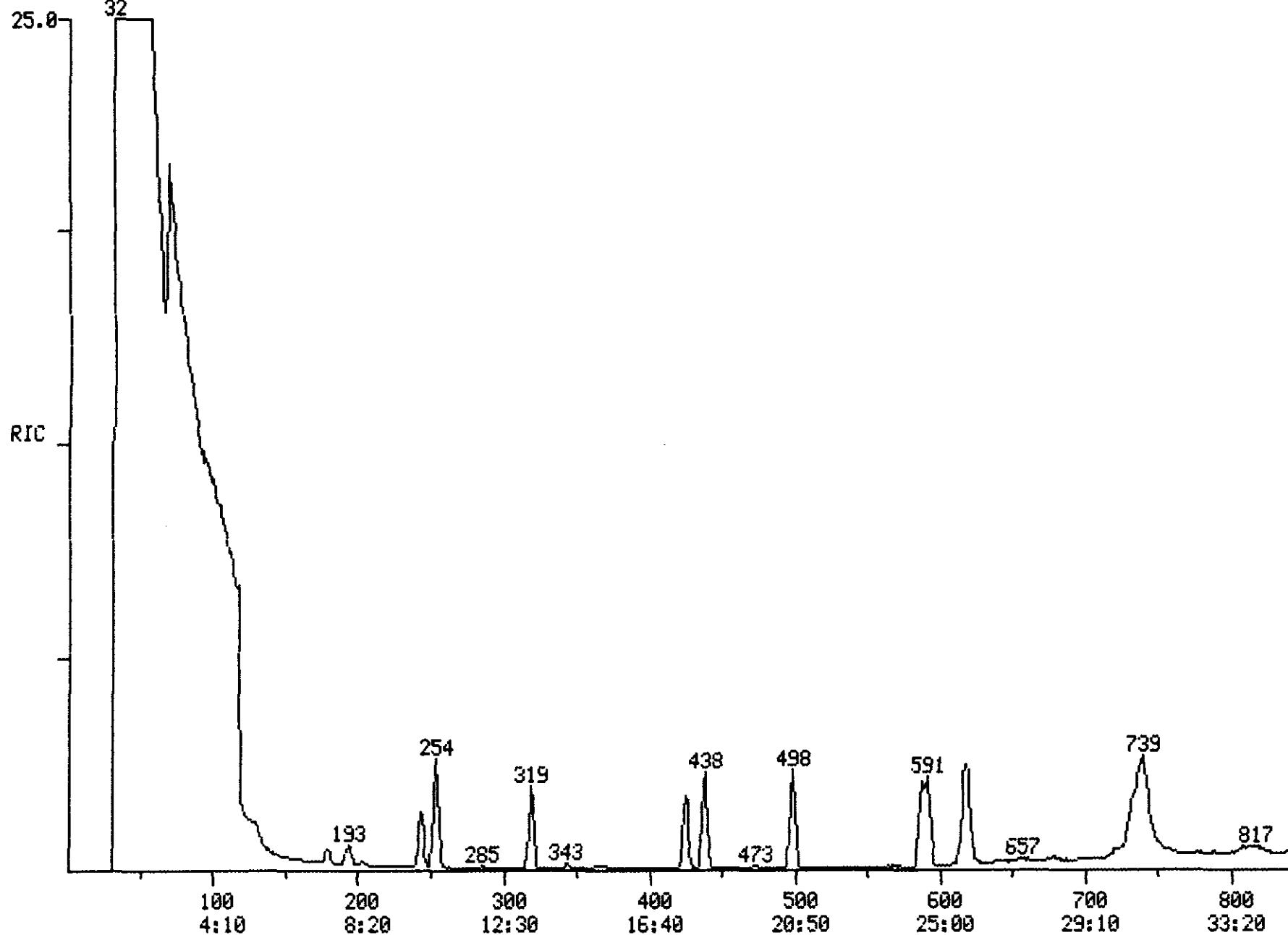
SAMPLE: CLP, VERSCOM, 2536, 1MSD, L, S, 16426MSD, U, , 420.1.0, 2,5G/5ML,

COND.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

RANGE: G 1, 840 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

100137

189440.



ORIGINATOR

(Req)

Quantitation Report File: U3399

Data: U3399.TI

04/26/90 18:53:00

Sample: CLP, VERSCDM, 2536, 1MSD, L, S, 16426MSD, V., 420.1.0, 2, 5G/5ML,

Conds.: INSTRUMENT U: SP-1000 COLUMN: 45C (3MIN) TO 225C @ 8C/MIN

Formula: -

Instrument: U

Weight: 0.011

Submitted by: VERSAR

Analyst: DD

Acct. No.: 420.1.0

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)

30.0 %

Resp. fac. from Library Entry

No	Name
1	C101 BROMOCHLOROMETHANE **IS#1**
2	C010 CHLOROMETHANE
3	C015 BROMOMETHANE
4	C020 VINYL CHLORIDE
5	C025 CHLOROETHANE
6	C030 METHYLENE CHLORIDE
7	C035 ACETONE
8	C040 CARBON DISULFIDE
9	C045 1,1-DICHLOROETHENE
10	C050 1,1-DICHLOROETHANE
11	C053 1,2-DICHLOROETHENE (TOTAL)
12	C060 CHLOROFORM
13	C065 1,2-DICHLOROETHANE
14	CS15 1,2-DICHLOROETHANE-D4 **SS#1**
15	C043 TRICHLOROFLUOROMETHANE
16	CI10 1,4-DIFLUOROBENZENE **IS#2**
17	C110 2-BUTANONE
18	C125 VINYL ACETATE
19	C120 CARBON TETRACHLORIDE
20	C130 BROMODICHLOROMETHANE
21	C140 1,2-DICHLOROPROPANE
22	C145 CIS-1,3-DICHLOROPROPENE
23	C150 TRICHLOROETHENE
24	C165 BENZENE
25	C155 DIBROMOCHLOROMETHANE
26	C160 1,1,2-TRICHLOROETHANE
27	C170 TRANS-1,3-DICHLOROPROPENE
28	C180 BROMOFORM
29	C115 1,1,1-TRICHLOROETHANE
30	CS05 TOLUENE-D8 **SS#2**
31	CI20 CHLOROBENZENE-D5 **IS#3**
32	C205 4-METHYL-2-PENTANONE
33	C210 2-HEXANONE
34	C220 TETRACHLOROETHENE
35	C225 1,1,2,2-TETRACHLOROETHANE
36	C230 TOLUENE
37	C235 CHLOROBENZENE
38	C240 ETHYLBENZENE
39	C245 STYRENE
40	C250 ORTHO & PARA XYLENE
41	C251 META-XYLENE
42	CS10 BROMOFLUOROBENZENE **SS#3**

all IS areas low
SS#1 recovery low
Ready

100138

ONLINE
(Raw)

U3399

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	128	253	10:32	1	1.000	A BB	18985.	50.000 UG/L*	8.92
2	NOT FOUND						low		
3	NOT FOUND								
4	NOT FOUND								
5	NOT FOUND								
6	84	180	7:30	1	0.711	A BB	4454.	8.550 UG/L	1.53
7	43	193	8:02	1	0.763	A BV	9447.	34.230 UG/L	6.11
8	NOT FOUND								
9	96	243	10:07	1	0.960	A BB	12589.	MS 28.788 UG/L	5.14
10	NOT FOUND								
11	NOT FOUND								
12	NOT FOUND								
13	NOT FOUND								
14	65	319	13:17	1	1.261	A BB	26826.	30.883 UG/L%	5.51 62
15	NOT FOUND								
16	114	498	20:45	16	1.000	A BB	46705.	50.000 UG/L*	8.92
17	NOT FOUND						low		
18	NOT FOUND								
19	NOT FOUND								
20	NOT FOUND								
21	NOT FOUND								
22	NOT FOUND								
23	130	425	17:42	16	0.853	A BB	15502.	MS 38.345 UG/L	6.84
24	78	438	18:15	16	0.880	A BB	46964.	MS 56.525 UG/L	10.09
25	NOT FOUND								
26	NOT FOUND								
27	NOT FOUND								
28	NOT FOUND								
29	NOT FOUND								
30	98	587	24:27	31	0.953	A BB	35072.	58.496 UG/L%	10.44 17
31	117	616	25:40	31	1.000	A BB	24935.	50.000 UG/L*	8.92
32	NOT FOUND						low		
33	NOT FOUND								
34	NOT FOUND								
35	NOT FOUND								
36	92	592	24:40	31	0.961	A BB	22088.	MS 64.532 UG/L	11.51
37	112	619	25:47	31	1.005	A BB	23561.	MS 51.655 UG/L	9.22
38	NOT FOUND								
39	NOT FOUND								
40	NOT FOUND								
41	NOT FOUND								
42	95	732	30:30	31	1.188	A BB	16685.	38.502 UG/L%	6.87 77

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
1	10:32	1.00	1.000	1.00	50.00	50.00	1.000	1.000	1.00
2	2:12		0.209						
3	3:22		0.320						
4	4:15		0.403						
5	5:15		0.498						
6	7:27	1.01	0.708	1.01	8.55	50.00	0.235	1.372	0.17
7	8:02	1.00	0.763	1.00	34.23	50.00	0.498	0.727	0.68
8	9:02		0.858						
9	10:07	1.00	0.960	1.00	28.79	50.00	0.663	1.152	0.58
10	11:25		1.083						
11	12:05		1.146						
12	12:40		1.202						

100139

U3399 ORIGINAL
(Red)

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
13	13:22		1. 269						
14	13:17	1. 00	1. 261	1. 00	30. 88	50. 00	1. 413	2. 288	0. 62
15	9:30		0. 901						
16	20:45	1. 00	1. 000	1. 00	50. 00	50. 00	1. 000	1. 000	1. 00
17	13:17		1. 261						
18	15:10		0. 731						
19	15:07		0. 729						
20	15:32		0. 749						
21	16:57		0. 817						
22	17:10		0. 827						
23	17:42	1. 00	0. 853	1. 00	38. 34	50. 00	0. 332	0. 433	0. 77
24	18:12	1. 00	0. 878	1. 00	56. 53	50. 00	1. 006	0. 889	1. 13
25	18:20		0. 884						
26	18:25		0. 888						
27	18:25		0. 888						
28	21:00		1. 012						
29	14:42		0. 709						
30	24:27	1. 00	0. 953	1. 00	58. 50	50. 00	1. 407	1. 202	1. 17
31	25:40	1. 00	1. 000	1. 00	50. 00	50. 00	1. 000	1. 000	1. 00
32	21:30		0. 838						
33	23:02		0. 898						
34	23:17		0. 907						
35	23:17		0. 907						
36	24:37	1. 00	0. 959	1. 00	64. 53	50. 00	0. 886	0. 686	1. 29
37	25:47	1. 00	1. 005	1. 00	51. 65	50. 00	0. 945	0. 915	1. 03
38	27:50		1. 084						
39	32:00		1. 247						
40	33:17		1. 297						
41	32:15		1. 256						
42	30:27	1. 00	1. 187	1. 00	38. 50	50. 00	0. 669	0. 869	0. 77

100140

VI. SAMPLE PREPARATION PACKAGE

1. Parameter Request Sheet
2. Screening Data Summary (soils only)
3. Comments
 - a. Extraction
 - b. Volatiles
 - c. Semivolatiles
 - d. Pesticides
4. Extraction Worksheets
 - a. Chronicle
 - b. Dry Weight Factor (soils only)
 - c. Dilution Factor (B) Worksheets
5. Injection Sequence Logs
6. Chain of Custody

4/29

VERSAR LABORATORY OPERATIONS
*** REVISED PARAMETER REQUEST SHEET ***

Deliver To: 48

Control: 2536
Code: VERSCDM
Batch: 2
Job Number: 420.1.0

Case: R3-7
SDG: 1
Quote: 9000194 - 90
Charge:

Manager: BECKMAN
Company: VERSAR DIV 31

Site: CDM

Received: 19-APR-90
Due: 17-MAY-90

Sampled: 18-APR-90
Report: 17-MAY-90

Comments: Protocols-Organic SOW 2/88; Inorganic Metals SOW 787 and CN SOW 788; TCLP by the new reg. Do sample QC on field # 1.

Test	Sample X C	Field Number	Matrix	Location	Other Tests
BNAL - 48					
	16419	A 1	SLUDGE	C-82	XBNA
	16420	A 2	SLUDGE	C-82	XBNA
	16421	A 3	SLUDGE	C-82	XBNA
	16422	A 5	SLUDGE	C-82	XBNA
	16423	B 6	SLUDGE	C-82	XBNA
	16424	B 7	SLUDGE	C-82	XBNA
CLPM - 43					
	16401	A 1	SLUDGE	C-82	
	16402	A 2	SLUDGE	C-82	
	16403	A 3	SLUDGE	C-82	
	16404	A 5	SLUDGE	C-82	
	16405	B 6	SLUDGE	C-82	
	16432	B 7	SLUDGE	C-82	
CNM - 44					
	16406	A 1	SLUDGE	C-82	TS
	16407	A 2	SLUDGE	C-82	TS
	16408	A 3	SLUDGE	C-82	TS
	16409	A 5	SLUDGE	C-82	TS
	16410	B 6	SLUDGE	C-82	TS
	16411	B 7	SLUDGE	C-82	TS
PEST - 47					
	16412	A 1	SLUDGE	C-82	XPES
	16413	A 2	SLUDGE	C-82	XPES
	16414	A 3	SLUDGE	C-82	XPES
	16415	A 5	SLUDGE	C-82	XPES
	16416	B 6	SLUDGE	C-82	XPES
	16417	B 7	SLUDGE	C-82	XPES
TCBN - 48					
	16418	B 1	SLUDGE	C-82	TCLM TCPs XTCH
TCHB - 47					
	16418	B 1	SLUDGE	C-82	TCLM XTCB XTCH
TCLM - 43					

07/13/86
(Ran)

Control #: 2536

Test	Sample	X	C	Field Number	Matrix	Location	Other Tests		
	16418	B	1		SLUDGE	C-82	TCLX	TCBN	XTCB
							TCPS	XTCP	TCHB
TCLV - 48	16425	B	1		SLUDGE	GCMS	ZHE		
TCLX - 43	16418	B	1		SLUDGE	C-82	TCLM	TCBN	XTCB
							TCPS	XTCP	TCHB
TCPS - 47	16418	B	1		SLUDGE	C-82	TCLM	TCLX	TCBN
							XTCB	XTCP	TCHB
TS - 44	16406	A	1		SLUDGE	C-82	CNM		
	16407	A	2		SLUDGE	C-82	CNM		
	16408	A	3		SLUDGE	C-82	CNM		
	16409	A	5		SLUDGE	C-82	CNM		
	16410	B	6		SLUDGE	C-82	CNM		
	16411	B	7		SLUDGE	C-82	CNM		
VOAL - 48	16426	B	1		SLUDGE	GCMS			
	16427	B	2		SLUDGE	GCMS			
	16428	B	3		SLUDGE	GCMS			
	16429	B	5		SLUDGE	GCMS			
	16430	B	6		SLUDGE	GCMS			
	16431	B	7		SLUDGE	GCMS			
XBNA - 46	16419	A	1		SLUDGE	C-82	BNAL		
	16420	A	2		SLUDGE	C-82	BNAL		
	16421	A	3		SLUDGE	C-82	BNAL		
	16422	A	5		SLUDGE	C-82	BNAL		
	16423	B	6		SLUDGE	C-82	BNAL		
	16424	B	7		SLUDGE	C-82	BNAL		
XPES - 46	16412	A	1		SLUDGE	C-82	PEST		
	16413	A	2		SLUDGE	C-82	PEST		
	16414	A	3		SLUDGE	C-82	PEST		
	16415	A	5		SLUDGE	C-82	PEST		
	16416	B	6		SLUDGE	C-82	PEST		
	16417	B	7		SLUDGE	C-82	PEST		
XTCB - 46	16418	B	1		SLUDGE	C-82	TCLM	TCLX	TCBN
							TCPS	XTCP	TCHB
XTCH - 46	16418	B	1		SLUDGE	C-82	TCLM	TCLX	TCBN
							XTCB	TCPS	XTCP
XTCP - 46	16418	B	1		SLUDGE	C-82	TCLM	TCLX	TCBN
							XTCB	TCPS	TCHB
ZHE - 43	16425	B	1		SLUDGE	GCMS	TCLV		

Versar[®] Inc.

Date: 5/1/90

TOTAL SOLIDS

Metler Balance Calibration

Zero	0.0000	0.0000
50 mg wt.	0.0500	0.0500
1 gram wt.	1.0000	1.0000
50 gram wt.	49.9989	49.9989

Project No. :
420.1-2 # 2536

pages: _____

NBS
Standard Ref. Material
1632b
T.V. = 98.7%

100144

mf
5/11/90

Analyst M. Ondersa
Date 5/8/90

VERSAR, INC.

GCMS INSTRUMENT ID 1
DATE 4-23-90

INJECTION LOG

Sectors Left 8394

Filename	ASCII Filename	Stream ID	Method	Case	SDG	Field Sample ID	Level	Matrix	Versar Sample No.	Frac.	Type	Project	Batch	Dil. F/T	Dil. F/T	Dil. F/T	Ext. Vol.	Purged	Purged	Final Dil.	IS No.
U3311		130				BFB 50ng														2ul	
U3312		130				BFB 50ng	tuned													2ul	
U3313		130	CLP			VSTD 50	L	S	22517	V										5ml	22388
U3314		130				BFB 50ng														2ul	
U3315		130				BFB 50ng														2ul	
U3316		130				BFB 50ng														2ul	
U3317		130				BFB 50ng														2ul	
U3318		130				BFB 50ng														2ul	
U3319		130				BFB 50ng														2ul	
U3320		130				BFB 50ng														2ul	
U3321		130				BFB 50ng														2ul	
U3322		130				BFB 50ng														2ul	
U3323		130				BFB 50ng														2ul	
U3324		130				BFB 50ng														2ul	
U3325		130				BFB 50ng														2ul	
U3326	T	130				BFB 50ng														2ul	

tuned @ 3:49 pm

Analyst Signature

D. Duse

Date 4-23-90

Original
CopyC
O
D
E
5

0000044

VERSAR, INC.

GCMS INSTRUMENT ID

DATE 4-23-90

INJECTION LOG

4042390-91

Filename	ASCII Filename	Stream ID	Method	Case	SDG	Field Sample ID	Level	Matrix	Versar Sample No.	Frac.	Type	Project	Batch	Ofl. F/T	Ofl. F/T	Ofl. F/T	Ext.	Vol. Purged	Final Ofl.	IS No.
U3322		130	CUP			NSTD50	L	S	22512	V	100						5ml		22385	
U3328		130	CUP			NSTD 200	L	S	22523	V	100						5ml		121	
U3329		130	CUP			NSTD 150	L	S	22524	V	100						5ml		1	
U3330		130	CUP			NSTD 100	L	S	22525	V	100						5ml			
U3331		130	CUP			NSTD20	L	S	22526	V	100						5ml			
U3332		130	CUP			NBLK32	L	S	NBLK32	V	SLR						5ml		22385	
U3333		130	CUP	VERS NOSH 2533		48442	L	S	16330A	V	-	420. 1.0					5ml		1	
U3334		130	CUP	PRC CEDR 2477		KS601001	L	S	14948	V	-	420. 61.0					18/5ml	1/5		
U3335		130	CUP	VERS NOSH 2533		48442 RE	L	S	16330A	V	-	420. 1.0					5g/5ml			

Analyst Signature

Alma

Date 4-23-90

000055

VERSAR, INC.

GCMS INSTRUMENT ID 4
DATE 4-26-90

INJECTION LOG

Sectors Left 7915

Filename	ASCII Filename	Stream ID	Method	Case	SDG	Field Sample ID	Level	Matrix	Versar Sample No.	Frac.	Type	Project	Batch	Dil. F/T	Dil. F/T	Dil. F/T	Ext.	Vol. Purged	Final Dil.	IS No.
3 U3385	T	130	CLP			BFB 50ng		tuned	@			7:52						2ul		
+ U3386	STD	130	CLP			VSTD 50	L	S	22546	V								5ml	22	
+ U3387	BLANK	130	CLP			VBLK 87	L	S	VBLK87	V								5ml	223	
U3388		130	CLP	ENGI HARL	2542	H V-SED-1	L	S	16627	V	-	420. 52.0	5				5g /5ml	223		
U3389		130	CLP	ENGI HARL	2542	HV-SED-2	L	S	16628	V	-	420. 52.0	5				5g /5ml	223		
U3390		130	CLP	ENGI ALBA	2544	AL-GW-5D-SI	L	S	16641	V	-	420. 52.0	3				5g /5ml	223		
U3391		130	CLP	ENGI ALBA	2562	AL-GW-40-SI	L	S	16895	V	-	420. 52.0	4				5g /5ml	223		
U3392		130	CLP	VERS CDM	2536	B 1	L	S	16426	V	-	420.1. 0	2				5g /5ml	223		
U3393		130	CLP	VERS CDM	2536	B 2	L	S	16427	V	-	420.1. 0	2				5g /5ml	223		
U3394		130	CLP	VERS CDM	2536	B 3	L	S	16428	V	-	420.1. 0	2				5g /5ml	223		
NB U3395		130	CLP	VERS CDM	2536	5 *	L	S	16429	V	-	420.1. 0	2				5g /5ml	223		
NB U3396		130	CLP	VERS CDM	2536	6	L	S	16430	V	-	420.1. 0	2				5g /5ml	223		
U3397		130	CLP	VERS CDM	2536	7	L	S	16431	V	-	420.1. 0	2				5g /5ml	223		
U3398	ms	130	CLP	VERS CDM	2536	1 MS	L	S	16426	V	-	420.1. 0	2				5g /5ml	223		
U3399	msd	130	CLP	VERS CDM	2536	1 MSD	L	S	16426	V	-	420.1. 0	2				5g /5ml	223		
ALST U3400		130	CLP	E+E AVV/E	2405	EWWP 372	L	S	177100	V	-	420. 51.0	96	100 ul / 5 ml			5ml	223		

30

(Instrument RESET)

Analyst Signature

Drew

Date 4-26-90

00147

VERSAR, INC.

GCMS INSTRUMENT ID U
DATE 4-26-90
MAT

INJECTION LOG

Sectors F4700

Filename	ASCII Filename	Stream ID	Method	Case	SDG	Field Sample ID	Level	Matrix	Versar Sample No.	Frac.	Type	Project	Batch	Dil. F/T	Dil. F/T	Dil. F/T	Ext. Purged	Vol. Purged	Final Dil.	IS No.
U3410	T					BFB 50ng		tuned	(2)			7.45						2ml		
U3411	MAT 30 4/67	130	CLP			MFT VS ID 50	L	S	22554	V	-	420.	1.0	2				5ml	22388	
U3412	BLK	130	CLP			VBLK12	L	S	VBLK12	V	BLK							5ml	22388	
U3413		130	CLP	VERS CDM	2536	B5	L	S	16429	V	-	420.	1.0	2				5g	22388	
U3414		130	CLP	VERS CDM	2536	B6	L	S	16430	V	-	420.	1.0	2				5g	22388	
U3415		130	CLP	VERS CDM	2536	B7	L	S	16431	V	-	420.	1.0	2				5g 5ml	22388	
U3416		130	CLP	GTI YORK	2453	SSI-3B	L	S	14579	V	-	420.	33.0	2				5g 5ml	22388	
U3417		130	CLP	GTI YORK	2453	SSI-4B	L	S	14580	V	-	420.	33.0	2				5g 5ml	22388	
U3418		130	CLP	GTI YORK	2454	SSI-6B	L	S	14606	V	-	420.	33.0	3				5g 5ml	22388	
U3419		130	CLP	GTI YORK	2454	SSI-10B	L	S	14608	V	-	420.	33.0	3				5g 5ml	22388	
U3420	**	130	CLP	VERS	2536	1	L	S	16426	V	-	420.	1.0	2				5g 5ml	22388	
U3421		130	CLP	ICP NEW	2549	SS29	L	S	16723	V	-	420.	58.0	10				5g/5ml	22388	
U3422		130	CLP			SS30	L	S	16724	V	-							5g/5ml	22388	
U3423		130	CLP			SS31	L	S	16725	V	-							5g/5ml	22388	
U3424		130	CLP			SS29 MS	L	S	16723MS	V	-							5g/5ml	22388	
U3425		130	CLP			SS29 MSD	L	S	16723MSD	V	-							5g/5ml	22388	

Relinquished by: (Signature)	Date / Time	Received by: (Signature)	Relinquished by: (Signature)	Date / Time	Received by: (Signature)
Sharon E. Schaeffer	4/18/90 2000				
Relinquished by: (Signature)	Date / Time	Received by: (Signature)	Relinquished by: (Signature)	Date / Time	Received by: (Signature)
100-1		R.D. Miron	4/19/90 0900	Slipped via Federal Express Airbill # 2294123705	

Distribution: Original Accompanies Shipment; Copy to Coordinator, Field Files

Remarks
Shipped via Federal Express
Airbill # 22941 23705

34

3-18549

HAIN OF CUSTODY RECORD

CHAIN OF CUSTODY RECORD

Relinquished by: (Signature)

Date / Time

Received by: (Signature)

Relinquished by: *(Signature)*

Date / Time

Received by: (Signature)

Sharon E. Schaefer

41870 2000

100

10. The following table summarizes the results of the study.

Relinquished by: (Signature)

Date / Time

Received by: (Signature)

Relinquished by: (Signature)

Date / Tim

Received by: (Signature)

BELINQUISHER BY: (Signature)

Date / Time

Received for Laboratory by
(Signature)

Date / Time
4/19/91 0900

Remark

Shipped via Federal Express
Airbill # 4024936093

CHAIN OF CUSTODY RECORD

CHAIN OF CUSTODY RECORD

PROJ. NO.		PROJECT NAME		NO. OF CON- TAINERS							REMARKS				
R3-7		R3-7			Sample Type										
SAMPLERS: (Signature)		Sharon E. Schaeffer, Lorraine B. Wyle			VOC		PCB/PCST.		Metals						
					TBVA		TALL		Chromate						
STA. NO.	DATE	TIME	COMP.	GRAB	STATION LOCATION										
2	4/18/90	11:35	X		Location 2						4	8 1 1 1 1	3-1068359, 63, 44, 65, 66, 67		
3	4/18/90	11:45	X		Location 3						4	2 1 1 1 1	3-1068368, 69, 70, 71, 72, 73		
5	4/18/90	14:30	X		Location 5						4	2 1 1 1 1	3-1068374, 75, 76, 77, 78, 79		
6	4/18/90	14:55	X		Location 6						5/18/90 10:10	2 1 2 2 2	3-1068380, 81, 82, 83, 84, 85, 86, 87, 88, 89		
7	4/18/90	15:05	X		Location 7						5/18/90 10:10	2 1 2 2 2	3-1068390, 91, 92, 93, 94, 95, 96, 97, 98, 99		
Relinquished by: (Signature)				Date / Time		Received by: (Signature)		Relinquished by: (Signature)				Date / Time		Received by: (Signature)	
Sharon E. Schaeffer				4/18/90 2000											
Relinquished by: (Signature)				Date / Time		Received by: (Signature)		Relinquished by: (Signature)				Date / Time		Received by: (Signature)	
						J. Z. Morrison									
Relinquished by: (Signature)				Date / Time		Received for Laboratory by: (Signature)		Date / Time		Remarks					
						J. Z. Morrison		4/19/90 0900							

CHAIN OF CUSTODY RECORD

Relinquished by: (Signature)

Date / Time

Received by: (Signature)

Being relinquished by: (Signature)

Date / Time

Received by: (Signature)

100-1000 mg/day.

- 7 -

Digitized by srujanika@gmail.com

Digitized by srujanika@gmail.com

— 1 —

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Relinquished by: (Signature)

Date / Time

Received for Laboratory by:

Page 1 Date / June

Remarks

Distribution: Original Accompanies Shipment, Copy to Coordinator, Field Files

Remarks
Shipped via Federal Express
Airbill # 2294123705